Full wwPDB X-ray Structure Validation Report (i)

Sep 25, 2023 – 07:28 AM EDT

PDB ID	:	5VU6
Title	:	TNA polymerase binary complex with primer/template duplex
Authors	:	Chim, N.; Chaput, J.C.
Deposited on	:	2017-05-18
Resolution	:	3.00 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
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) 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 (i)) were used in the production of this report:

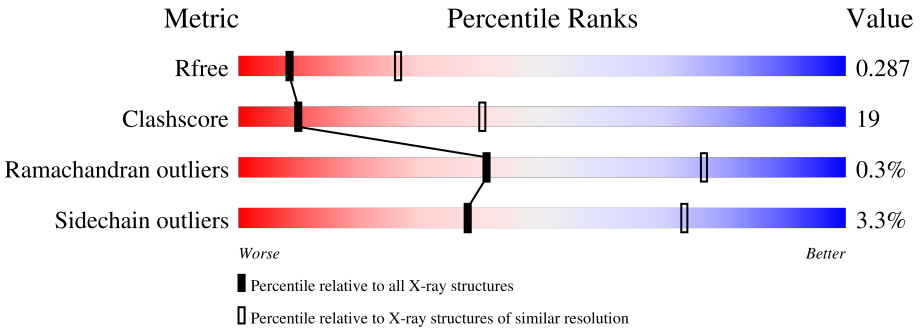
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

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 3.00 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}	130704	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)

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 <=5%

Mol	Chain	Length	Quality of chain
1	A	774	62% 33% • •
2	T	16	38% 50% 12%
3	P	12	50% 42% 8%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 6743 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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	756	6214	4003	1053	1142	16	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	141	ALA	ASP	engineered mutation	UNP D0VWU9
A	143	ALA	GLU	engineered mutation	UNP D0VWU9
A	147	HIS	GLU	engineered mutation	UNP D0VWU9
A	485	ARG	ALA	engineered mutation	UNP D0VWU9
A	584	LYS	GLU	engineered mutation	UNP D0VWU9
A	664	ILE	GLU	engineered mutation	UNP D0VWU9

- Molecule 2 is a DNA chain called DNA template.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	T	14	286	136	50	86	14	0	0	0

- Molecule 3 is a DNA chain called DNA/TNA hybrid primer.

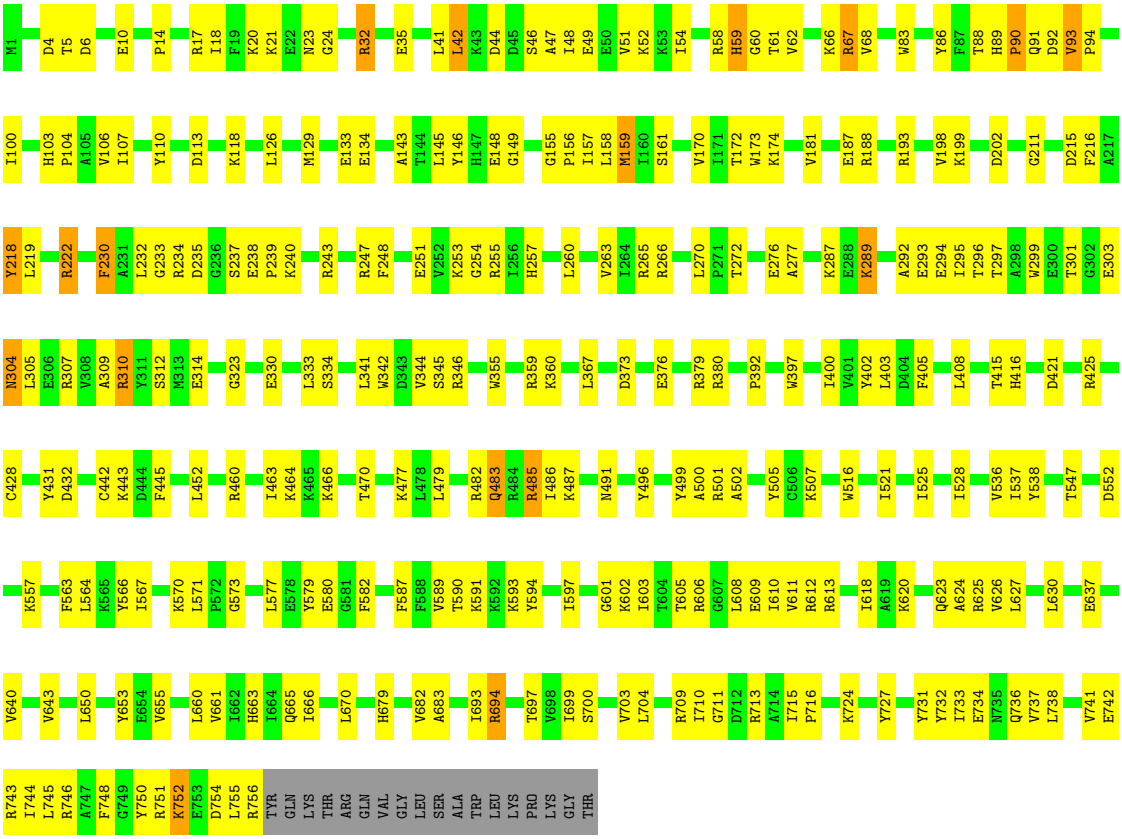
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
3	P	12	243	115	49	68	11	0	0	0

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. 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. 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

Chain A: 62% 33%

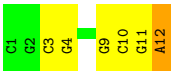


- Molecule 2: DNA template

Chain T: 38% 50% 12%

- Molecule 3: DNA/TNA hybrid primer

Chain P: 50% 42% 8%


C1
G2
C3
G4
G9
C10
G11
A12

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, α, β, γ	66.09Å 110.80Å 149.08Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	88.93 – 3.00 88.93 – 2.59	Depositor EDS
% Data completeness (in resolution range)	99.7 (88.93-3.00) 79.2 (88.93-2.59)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	0.19 (at 2.58Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, R_{free}	0.235 , 0.283 0.250 , 0.287	Depositor DCC
R_{free} test set	2000 reflections (5.79%)	wwPDB-VP
Wilson B-factor (Å^2)	37.4	Xtriage
Anisotropy	0.520	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.39 , 50.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.88	EDS
Total number of atoms	6743	wwPDB-VP
Average B, all atoms (Å^2)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.66% 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, 0.333 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: FA2

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	2/6353 (0.0%)	0.53	2/8570 (0.0%)
2	T	0.58	0/319	0.90	0/490
3	P	0.62	0/250	0.86	0/384
All	All	0.42	2/6922 (0.0%)	0.57	2/9444 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	42	LEU	C-N	9.48	1.55	1.34
1	A	218	TYR	CE1-CZ	-5.37	1.31	1.38

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	42	LEU	O-C-N	-6.38	112.49	122.70
1	A	304	ASN	N-CA-C	-5.67	95.70	111.00

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	A	6214	0	6264	244	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	T	286	0	159	10	0
3	P	243	0	132	8	0
All	All	6743	0	6555	250	1

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 19.

All (250) 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:A:610:ILE:HD12	1:A:624:ALA:HB2	1.42	1.01
1:A:47:ALA:O	1:A:51:VAL:HG23	1.74	0.88
1:A:303:GLU:HG2	1:A:304:ASN:H	1.40	0.87
1:A:158:LEU:HD23	1:A:299:TRP:CE3	2.20	0.77
1:A:88:THR:OG1	1:A:92:ASP:OD2	2.01	0.75
1:A:49:GLU:OE1	1:A:52:LYS:HD3	1.87	0.75
1:A:42:LEU:HD13	1:A:44:ASP:O	1.85	0.74
1:A:145:LEU:HD23	1:A:156:PRO:HD2	1.70	0.73
1:A:738:LEU:O	1:A:742:GLU:N	2.20	0.73
1:A:405:PHE:HD2	1:A:408:LEU:HD13	1.51	0.73
1:A:609:GLU:OE2	1:A:743:ARG:NH1	2.21	0.73
1:A:567:ILE:HD11	1:A:577:LEU:HB3	1.71	0.72
1:A:211:GLY:HA2	1:A:215:ASP:HB2	1.72	0.72
1:A:93:VAL:HB	1:A:94:PRO:HD3	1.72	0.72
1:A:155:GLY:H	1:A:222:ARG:HH22	1.37	0.72
1:A:303:GLU:O	1:A:304:ASN:C	2.28	0.71
1:A:23:ASN:N	1:A:133:GLU:OE1	2.24	0.71
1:A:303:GLU:O	1:A:305:LEU:N	2.24	0.70
1:A:643:VAL:HG11	1:A:755:LEU:HD13	1.74	0.70
1:A:587:PHE:CE1	1:A:744:ILE:HG23	2.26	0.70
1:A:460:ARG:HG3	1:A:483:GLN:HG2	1.73	0.69
1:A:587:PHE:CZ	1:A:744:ILE:HG23	2.28	0.69
1:A:21:LYS:HE3	1:A:24:GLY:HA2	1.75	0.69
1:A:126:LEU:O	1:A:359:ARG:NH2	2.26	0.68
1:A:650:LEU:HD22	1:A:655:VAL:HG21	1.76	0.68
1:A:199:LYS:NZ	1:A:233:GLY:O	2.27	0.68
1:A:589:VAL:HG22	1:A:744:ILE:HG12	1.76	0.68
1:A:611:VAL:HG12	3:P:10:DC:H5"	1.75	0.67
1:A:376:GLU:HA	1:A:379:ARG:HD3	1.76	0.67
1:A:161:SER:HG	1:A:312:SER:HG	1.39	0.67
1:A:610:ILE:CD1	1:A:624:ALA:HB2	2.20	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:733:ILE:HA	1:A:737:VAL:HG12	1.77	0.66
1:A:270:LEU:O	1:A:613:ARG:NH1	2.28	0.66
1:A:710:ILE:HD12	2:T:12:DT:H3'	1.78	0.65
1:A:129:MET:HG3	1:A:334:SER:OG	1.97	0.64
1:A:743:ARG:NH2	2:T:11:DT:OP1	2.30	0.64
1:A:392:PRO:HB3	1:A:538:TYR:HD1	1.62	0.64
1:A:52:LYS:HE3	1:A:68:VAL:HG21	1.79	0.64
1:A:59:HIS:C	1:A:61:THR:H	2.00	0.64
1:A:602:LYS:NZ	1:A:603:ILE:O	2.31	0.64
1:A:243:ARG:HG3	1:A:248:PHE:HE1	1.63	0.64
1:A:670:LEU:HD11	1:A:693:ILE:HB	1.79	0.64
1:A:303:GLU:HG2	1:A:304:ASN:N	2.12	0.63
1:A:263:VAL:CG2	1:A:345:SER:HB2	2.28	0.63
1:A:734:GLU:OE2	1:A:756:ARG:NH1	2.32	0.63
1:A:612:ARG:HD2	3:P:10:DC:H5'	1.81	0.63
1:A:145:LEU:HB2	1:A:158:LEU:HD11	1.79	0.62
1:A:742:GLU:HG3	1:A:752:LYS:HG3	1.80	0.62
1:A:297:THR:O	1:A:301:THR:HG22	2.00	0.62
1:A:59:HIS:O	1:A:61:THR:HG23	2.01	0.61
1:A:605:THR:HG23	1:A:608:LEU:HD12	1.83	0.61
1:A:157:ILE:N	1:A:187:GLU:OE1	2.34	0.61
1:A:257:HIS:O	1:A:346:ARG:NH2	2.33	0.61
1:A:670:LEU:HD12	1:A:670:LEU:H	1.66	0.60
1:A:487:LYS:O	1:A:491:ASN:ND2	2.32	0.60
2:T:5:DT:H3	3:P:12:FA2:H2	1.65	0.60
1:A:110:TYR:CD1	1:A:367:LEU:HD21	2.36	0.60
1:A:240:LYS:HB2	1:A:251:GLU:HB2	1.83	0.60
1:A:745:LEU:HA	1:A:748:PHE:HD2	1.66	0.60
1:A:605:THR:HG21	1:A:610:ILE:CG2	2.32	0.60
1:A:59:HIS:O	1:A:61:THR:N	2.34	0.60
1:A:397:TRP:HB2	1:A:400:ILE:HD11	1.84	0.60
1:A:143:ALA:HB3	1:A:159:MET:HG3	1.84	0.59
1:A:301:THR:HG23	1:A:303:GLU:H	1.68	0.59
1:A:650:LEU:CB	1:A:733:ILE:HD11	2.32	0.59
1:A:265:ARG:NH2	1:A:665:GLN:OE1	2.36	0.59
1:A:334:SER:HA	1:A:344:VAL:HG21	1.85	0.59
1:A:59:HIS:C	1:A:61:THR:N	2.56	0.58
1:A:380:ARG:NH1	1:A:500:ALA:O	2.32	0.58
1:A:663:HIS:CD2	1:A:700:SER:CB	2.86	0.58
1:A:6:ASP:OD2	1:A:253:LYS:NZ	2.36	0.58
1:A:610:ILE:HA	1:A:620:LYS:HG2	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:403:LEU:HD22	1:A:564:LEU:HD21	1.86	0.58
1:A:703:VAL:HG12	1:A:732:TYR:CE2	2.38	0.58
1:A:41:LEU:O	1:A:107:ILE:N	2.35	0.57
1:A:742:GLU:OE1	1:A:746:ARG:NH2	2.37	0.57
1:A:751:ARG:O	1:A:754:ASP:N	2.35	0.57
1:A:91:GLN:O	1:A:94:PRO:HD2	2.05	0.57
1:A:405:PHE:CD2	1:A:408:LEU:HD13	2.37	0.57
1:A:694:ARG:HG3	1:A:697:THR:HB	1.86	0.57
1:A:47:ALA:O	1:A:51:VAL:N	2.37	0.57
1:A:158:LEU:HD21	1:A:299:TRP:CG	2.40	0.56
1:A:243:ARG:HG3	1:A:248:PHE:CE1	2.40	0.56
1:A:660:LEU:O	1:A:732:TYR:OH	2.18	0.56
1:A:666:ILE:HG12	1:A:699:ILE:HD11	1.89	0.55
1:A:463:ILE:HG23	1:A:479:LEU:HD22	1.89	0.55
1:A:724:LYS:O	1:A:724:LYS:HG2	2.06	0.55
1:A:216:PHE:HE2	1:A:257:HIS:CD2	2.24	0.54
1:A:110:TYR:HD1	1:A:367:LEU:HD21	1.72	0.54
1:A:145:LEU:CD2	1:A:156:PRO:HD2	2.36	0.54
1:A:650:LEU:HB2	1:A:733:ILE:HD11	1.89	0.54
1:A:525:ILE:HD12	1:A:536:VAL:HG11	1.90	0.54
1:A:155:GLY:H	1:A:222:ARG:NH2	2.06	0.53
1:A:4:ASP:OD1	1:A:5:THR:N	2.37	0.53
1:A:466:LYS:O	1:A:470:THR:HG23	2.09	0.53
1:A:158:LEU:HD21	1:A:299:TRP:CB	2.39	0.53
1:A:158:LEU:CD2	1:A:299:TRP:CD2	2.92	0.53
1:A:590:THR:OG1	1:A:591:LYS:N	2.42	0.53
1:A:6:ASP:OD1	1:A:17:ARG:HB2	2.08	0.52
1:A:292:ALA:O	1:A:296:THR:HG22	2.09	0.52
1:A:155:GLY:N	1:A:222:ARG:HH22	2.06	0.52
1:A:416:HIS:ND1	1:A:432:ASP:OD1	2.38	0.52
1:A:18:ILE:HG22	1:A:20:LYS:HG3	1.91	0.52
1:A:303:GLU:CG	1:A:304:ASN:H	2.16	0.52
1:A:143:ALA:HB1	1:A:295:ILE:HD11	1.91	0.52
1:A:230:PHE:HB3	1:A:238:GLU:HG2	1.92	0.52
1:A:613:ARG:HG3	3:P:10:DC:OP1	2.10	0.52
1:A:89:HIS:CE1	1:A:91:GLN:HG3	2.45	0.52
1:A:234:ARG:HB2	1:A:254:GLY:HA3	1.92	0.52
1:A:464:LYS:HG3	1:A:483:GLN:HG3	1.92	0.52
1:A:129:MET:HE2	1:A:341:LEU:HD12	1.92	0.51
1:A:67:ARG:HG2	1:A:68:VAL:N	2.25	0.51
1:A:89:HIS:O	1:A:92:ASP:HB2	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:174:LYS:HE2	1:A:299:TRP:CE2	2.46	0.51
1:A:146:TYR:CE2	1:A:148:GLU:HA	2.45	0.51
1:A:159:MET:HE3	1:A:309:ALA:HA	1.93	0.51
1:A:663:HIS:CD2	1:A:700:SER:HB3	2.46	0.51
1:A:703:VAL:HG11	1:A:731:TYR:CD2	2.46	0.51
1:A:89:HIS:N	1:A:92:ASP:OD2	2.42	0.51
2:T:5:DT:H3	3:P:12:FA2:C2	2.23	0.51
1:A:158:LEU:HD23	1:A:299:TRP:CD2	2.46	0.51
1:A:263:VAL:HG21	1:A:345:SER:HB2	1.92	0.51
1:A:425:ARG:HB3	1:A:442:CYS:HB2	1.93	0.51
1:A:610:ILE:HA	1:A:620:LYS:HB3	1.93	0.51
1:A:103:HIS:O	1:A:106:VAL:HG12	2.12	0.50
1:A:593:LYS:HD2	2:T:9:DA:H4'	1.92	0.50
1:A:682:VAL:HG13	1:A:716:PRO:HD3	1.94	0.50
1:A:683:ALA:HB2	1:A:699:ILE:HD11	1.93	0.50
1:A:557:LYS:HE2	1:A:582:PHE:CD1	2.47	0.50
1:A:593:LYS:NZ	2:T:10:DG:OP1	2.44	0.50
1:A:597:ILE:HD13	1:A:603:ILE:HG12	1.94	0.50
1:A:521:ILE:O	1:A:525:ILE:HG12	2.11	0.50
1:A:35:GLU:OE1	1:A:66:LYS:NZ	2.45	0.50
1:A:466:LYS:HE2	1:A:479:LEU:HD11	1.94	0.50
1:A:653:TYR:HA	1:A:727:TYR:OH	2.11	0.49
1:A:41:LEU:HD12	1:A:42:LEU:H	1.76	0.49
1:A:743:ARG:HD3	2:T:10:DG:OP1	2.12	0.49
1:A:402:TYR:HD2	1:A:580:GLU:HB3	1.76	0.49
1:A:21:LYS:HE2	1:A:134:GLU:O	2.12	0.49
1:A:563:PHE:CZ	1:A:567:ILE:HG21	2.48	0.49
1:A:650:LEU:HB3	1:A:733:ILE:HD11	1.95	0.49
1:A:711:GLY:N	2:T:13:DC:OP1	2.42	0.49
1:A:158:LEU:CD2	1:A:299:TRP:CG	2.96	0.48
1:A:159:MET:HB3	1:A:172:THR:HB	1.95	0.48
1:A:421:ASP:OD1	1:A:421:ASP:N	2.45	0.48
1:A:44:ASP:C	1:A:46:SER:H	2.17	0.48
1:A:428:CYS:HB2	1:A:431:TYR:CZ	2.48	0.48
1:A:276:GLU:OE2	1:A:287:LYS:NZ	2.46	0.48
1:A:276:GLU:HG3	1:A:287:LYS:HG2	1.96	0.48
1:A:679:HIS:O	1:A:699:ILE:HD13	2.14	0.48
1:A:42:LEU:CD1	1:A:44:ASP:O	2.60	0.48
1:A:342:TRP:O	1:A:346:ARG:HG3	2.14	0.48
1:A:537:ILE:HD11	1:A:547:THR:HG22	1.95	0.48
1:A:403:LEU:HD23	1:A:579:TYR:HA	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:589:VAL:CG2	1:A:744:ILE:HG12	2.42	0.48
1:A:609:GLU:OE1	1:A:609:GLU:N	2.43	0.48
1:A:158:LEU:CD2	1:A:299:TRP:HB2	2.44	0.48
1:A:660:LEU:HB3	1:A:732:TYR:CZ	2.49	0.48
1:A:143:ALA:HB1	1:A:295:ILE:CD1	2.43	0.47
1:A:118:LYS:HD3	1:A:355:TRP:CH2	2.50	0.47
1:A:158:LEU:HD21	1:A:299:TRP:HB2	1.96	0.47
1:A:243:ARG:HA	1:A:248:PHE:CD1	2.49	0.47
1:A:683:ALA:HB2	1:A:699:ILE:CD1	2.44	0.47
1:A:155:GLY:O	1:A:218:TYR:OH	2.28	0.47
1:A:294:GLU:OE1	1:A:307:ARG:NH2	2.45	0.47
1:A:564:LEU:HD11	1:A:579:TYR:HB2	1.96	0.47
1:A:333:LEU:HB2	1:A:485:ARG:HD2	1.96	0.47
1:A:100:ILE:O	1:A:106:VAL:HG11	2.15	0.46
1:A:623:GLN:O	1:A:627:LEU:HB2	2.14	0.46
1:A:14:PRO:HG3	1:A:90:PRO:HD3	1.98	0.46
1:A:48:ILE:HG21	1:A:83:TRP:CZ2	2.50	0.46
1:A:272:THR:HG23	1:A:277:ALA:CB	2.46	0.46
1:A:477:LYS:HD2	1:A:477:LYS:HA	1.71	0.46
1:A:742:GLU:O	1:A:746:ARG:NE	2.48	0.46
1:A:505:TYR:OH	1:A:507:LYS:HB2	2.16	0.46
1:A:232:LEU:HB2	1:A:239:PRO:HD3	1.97	0.46
1:A:528:ILE:HB	1:A:563:PHE:CZ	2.51	0.46
1:A:58:ARG:O	1:A:59:HIS:ND1	2.41	0.46
1:A:392:PRO:HB3	1:A:538:TYR:CD1	2.48	0.46
1:A:640:VAL:HG11	1:A:750:TYR:CE1	2.51	0.46
1:A:222:ARG:HD3	1:A:222:ARG:HA	1.60	0.46
1:A:625:ARG:HA	1:A:625:ARG:HD3	1.67	0.45
1:A:737:VAL:O	1:A:741:VAL:HG22	2.16	0.45
1:A:145:LEU:HB2	1:A:158:LEU:CD1	2.46	0.45
1:A:376:GLU:O	1:A:379:ARG:NH1	2.44	0.45
1:A:432:ASP:HB2	1:A:443:LYS:HE3	1.34	0.45
1:A:113:ASP:N	1:A:113:ASP:OD1	2.48	0.45
1:A:709:ARG:O	1:A:713:ARG:HG3	2.15	0.45
1:A:704:LEU:HD11	1:A:715:ILE:HD13	1.99	0.44
2:T:5:DT:H2'	2:T:6:DC:C6	2.51	0.44
1:A:235:ASP:OD1	1:A:237:SER:OG	2.31	0.44
1:A:640:VAL:HG12	1:A:755:LEU:HD21	1.99	0.44
1:A:148:GLU:HG2	1:A:149:GLY:N	2.32	0.44
3:P:3:DC:H2"	3:P:4:DG:C8	2.53	0.44
1:A:202:ASP:HA	1:A:255:ARG:HH12	1.83	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:276:GLU:HG2	1:A:289:LYS:HB3	2.00	0.44
1:A:693:ILE:HG23	1:A:697:THR:HG21	1.99	0.44
1:A:42:LEU:C	1:A:42:LEU:HD12	2.38	0.44
1:A:496:TYR:CZ	1:A:502:ALA:HB1	2.52	0.44
1:A:627:LEU:CD1	1:A:744:ILE:HD13	2.48	0.44
1:A:597:ILE:HD11	1:A:601:GLY:HA2	1.99	0.43
1:A:626:VAL:O	1:A:630:LEU:HD23	2.18	0.43
1:A:216:PHE:CE2	1:A:257:HIS:CD2	3.05	0.43
1:A:360:LYS:HA	1:A:360:LYS:HD2	1.76	0.43
1:A:425:ARG:NE	1:A:445:PHE:CE1	2.87	0.43
1:A:174:LYS:HG3	1:A:299:TRP:CH2	2.53	0.43
1:A:170:VAL:HB	1:A:181:VAL:HG22	2.01	0.43
1:A:589:VAL:HG22	1:A:744:ILE:CG1	2.48	0.43
1:A:742:GLU:CD	1:A:746:ARG:HE	2.22	0.43
1:A:240:LYS:NZ	1:A:253:LYS:HE3	2.34	0.43
1:A:301:THR:HG23	1:A:303:GLU:N	2.34	0.43
1:A:589:VAL:HG13	1:A:743:ARG:O	2.19	0.43
1:A:159:MET:CE	1:A:309:ALA:HA	2.48	0.42
1:A:609:GLU:O	1:A:620:LYS:HG2	2.19	0.42
1:A:103:HIS:HD1	1:A:104:PRO:HD2	1.84	0.42
1:A:211:GLY:O	1:A:216:PHE:HB2	2.19	0.42
1:A:292:ALA:O	1:A:295:ILE:HG22	2.18	0.42
1:A:566:TYR:CE2	1:A:570:LYS:HD2	2.54	0.42
2:T:8:DC:H2"	2:T:9:DA:C8	2.54	0.42
1:A:289:LYS:HE2	1:A:289:LYS:HB2	1.80	0.42
1:A:630:LEU:HD23	1:A:630:LEU:H	1.84	0.42
1:A:732:TYR:HA	1:A:736:GLN:HB2	2.02	0.42
1:A:51:VAL:HA	1:A:54:ILE:HD11	2.01	0.42
1:A:310:ARG:NE	1:A:314:GLU:OE2	2.52	0.42
1:A:594:TYR:CZ	1:A:606:ARG:HG3	2.54	0.42
1:A:742:GLU:O	1:A:746:ARG:NH2	2.53	0.42
1:A:51:VAL:HG22	1:A:103:HIS:CD2	2.54	0.42
1:A:247:ARG:NH1	1:A:694:ARG:HD3	2.35	0.42
1:A:415:THR:HG23	1:A:573:GLY:HA3	2.01	0.42
1:A:663:HIS:CD2	1:A:700:SER:HB2	2.55	0.42
3:P:10:DC:H2'	3:P:11:DG:C8	2.55	0.42
1:A:10:GLU:OE2	1:A:32:ARG:NH1	2.53	0.41
1:A:661:VAL:CG1	1:A:700:SER:HB2	2.50	0.41
1:A:35:GLU:HB2	1:A:86:TYR:HD2	1.85	0.41
1:A:482:ARG:O	1:A:486:ILE:HG13	2.21	0.41
1:A:215:ASP:O	1:A:219:LEU:HD12	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:129:MET:CG	1:A:334:SER:OG	2.67	0.41
1:A:260:LEU:HD21	1:A:323:GLY:HA2	2.02	0.41
1:A:292:ALA:HA	1:A:295:ILE:HG22	2.02	0.41
1:A:360:LYS:HG3	1:A:452:LEU:HD22	2.02	0.41
1:A:571:LEU:HD23	1:A:571:LEU:HA	1.84	0.41
1:A:263:VAL:HG11	1:A:330:GLU:HG3	2.02	0.41
1:A:637:GLU:H	1:A:637:GLU:HG2	1.66	0.41
1:A:499:TYR:OH	1:A:501:ARG:HD3	2.20	0.41
1:A:612:ARG:NH1	3:P:9:DG:H1'	2.35	0.41
1:A:158:LEU:C	1:A:159:MET:HG2	2.40	0.41
1:A:198:VAL:O	1:A:202:ASP:N	2.46	0.40
1:A:48:ILE:O	1:A:52:LYS:N	2.55	0.40
1:A:89:HIS:HE1	1:A:91:GLN:HG3	1.85	0.40
1:A:251:GLU:HG3	1:A:342:TRP:CZ2	2.57	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:293:GLU:CB	1:A:724:LYS:NZ[2_557]	2.04	0.16

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	754/774 (97%)	720 (96%)	32 (4%)	2 (0%)	41 76

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	60	GLY
1	A	93	VAL

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
1	A	658/673 (98%)	636 (97%)	22 (3%)	38 73

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	ARG
1	A	59	HIS
1	A	62	VAL
1	A	67	ARG
1	A	90	PRO
1	A	159	MET
1	A	173	TRP
1	A	188	ARG
1	A	193	ARG
1	A	222	ARG
1	A	230	PHE
1	A	266	ARG
1	A	289	LYS
1	A	310	ARG
1	A	373	ASP
1	A	483	GLN
1	A	485	ARG
1	A	516	TRP
1	A	552	ASP
1	A	618	ILE
1	A	694	ARG
1	A	752	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	663	HIS

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	$\#\|Z\| > 2$
3	FA2	P	12	2,3	16,22,23	6.16	12 (75%)	15,32,35	3.12	6 (40%)

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
3	FA2	P	12	2,3	-	1/1/21/22	0/3/3/3

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	P	12	FA2	C2-N3	13.26	1.53	1.32
3	P	12	FA2	C4-N3	10.39	1.50	1.35
3	P	12	FA2	O2'-C2'	-10.07	1.19	1.43
3	P	12	FA2	C6-C5	-7.00	1.17	1.43
3	P	12	FA2	C8-N7	-6.73	1.22	1.34
3	P	12	FA2	C6-N1	-5.14	1.14	1.37
3	P	12	FA2	C5-C4	4.56	1.53	1.40
3	P	12	FA2	O4'-C1'	4.49	1.48	1.41
3	P	12	FA2	C2'-C1'	-4.17	1.47	1.53
3	P	12	FA2	O4'-C4'	3.89	1.52	1.43
3	P	12	FA2	C2'-C3'	-3.44	1.44	1.53
3	P	12	FA2	C6-N6	3.24	1.45	1.34

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	P	12	FA2	C4-C5-N7	-8.41	100.64	109.40
3	P	12	FA2	C2-N1-C6	5.41	128.01	118.75
3	P	12	FA2	N3-C2-N1	-3.31	123.50	128.68
3	P	12	FA2	C1'-N9-C4	-3.17	121.08	126.64
3	P	12	FA2	C4'-C3'-C2'	3.08	107.66	102.28
3	P	12	FA2	C3'-C2'-C1'	2.84	106.22	99.92

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	P	12	FA2	C2'-C3'-O3'-P

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	P	12	FA2	2	0

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

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 (i)

6.1 Protein, DNA and RNA chains (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers (i)

Unable to reproduce the depositors R factor - this section is therefore empty.