



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:30 PM GMT

PDB ID : 4LT6
Title : Crystal Structure of human poly(A) polymerase gamma
Authors : Yang, Q.; Nausch, L.; Martin, G.; Keller, W.; Doublié, S.
Deposited on : 2013-07-23
Resolution : 2.79 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (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

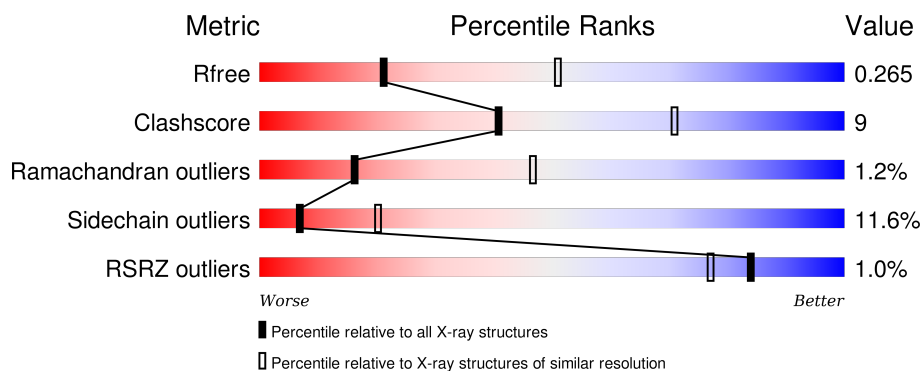
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.79 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	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

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 ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	514	
1	B	514	

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
2	3AT	A	601	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7644 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.

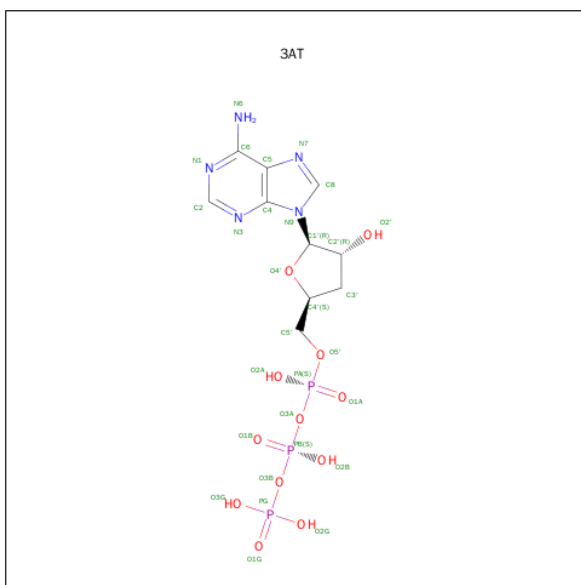
- Molecule 1 is a protein called Poly(A) polymerase gamma.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	467	Total	C	N	O	S	0	0	0
			3784	2436	652	684	12			
1	B	462	Total	C	N	O	S	0	0	0
			3745	2410	646	677	12			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	HIS	-	EXPRESSION TAG	UNP Q9BWT3
A	-4	HIS	-	EXPRESSION TAG	UNP Q9BWT3
A	-3	HIS	-	EXPRESSION TAG	UNP Q9BWT3
A	-2	HIS	-	EXPRESSION TAG	UNP Q9BWT3
A	-1	HIS	-	EXPRESSION TAG	UNP Q9BWT3
A	0	HIS	-	EXPRESSION TAG	UNP Q9BWT3
B	-5	HIS	-	EXPRESSION TAG	UNP Q9BWT3
B	-4	HIS	-	EXPRESSION TAG	UNP Q9BWT3
B	-3	HIS	-	EXPRESSION TAG	UNP Q9BWT3
B	-2	HIS	-	EXPRESSION TAG	UNP Q9BWT3
B	-1	HIS	-	EXPRESSION TAG	UNP Q9BWT3
B	0	HIS	-	EXPRESSION TAG	UNP Q9BWT3

- Molecule 2 is 3'-DEOXYADENOSINE-5'-TRIPHOSPHATE (three-letter code: 3AT) (formula: C₁₀H₁₆N₅O₁₂P₃).



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

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Ca	0	0
			1	1		
3	A	1	Total	Ca	0	0
			1	1		

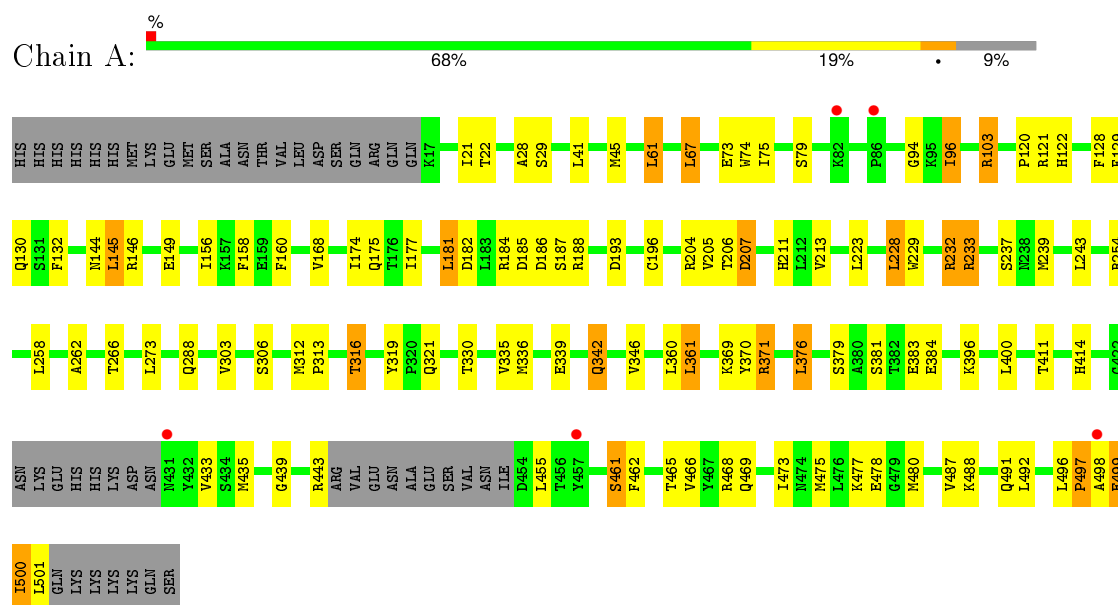
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	27	Total	O	0	0
			27	27		
4	B	26	Total	O	0	0
			26	26		

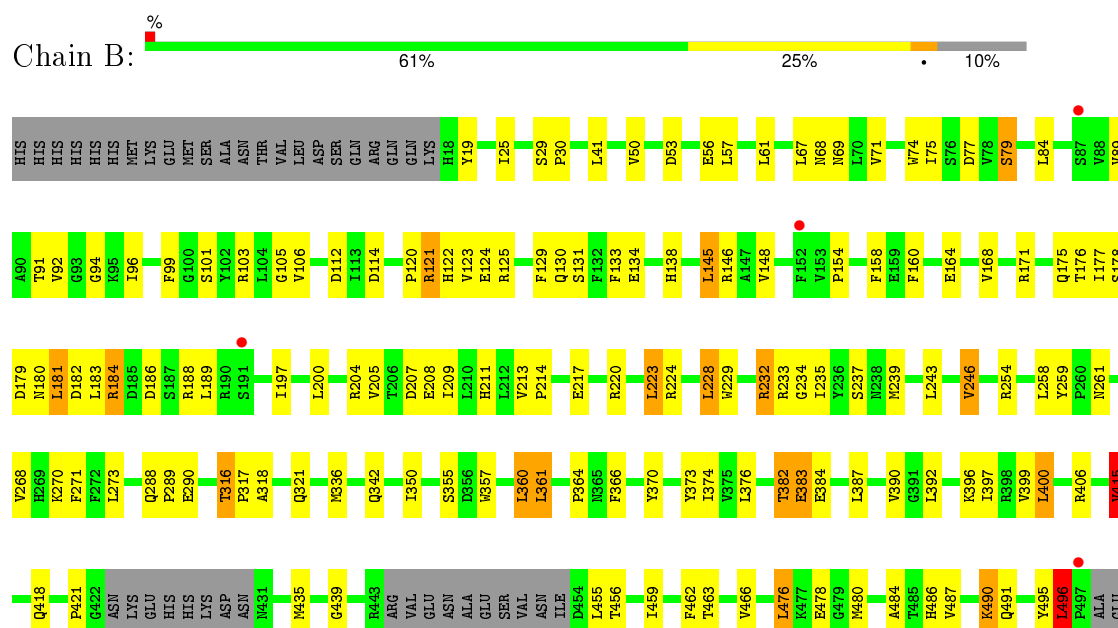
3 Residue-property plots

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: Poly(A) polymerase gamma



- Molecule 1: Poly(A) polymerase gamma



ILE
LEU
GLN
LYS
LYS
LYS
LYS
GLN
SER

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	68.87Å 89.92Å 201.63Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.88 – 2.79 19.84 – 2.79	Depositor EDS
% Data completeness (in resolution range)	95.7 (19.88-2.79) 89.9 (19.84-2.79)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.65 (at 2.79Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.193 , 0.257 0.203 , 0.265	Depositor DCC
R_{free} test set	1507 reflections (5.38%)	DCC
Wilson B-factor (Å ²)	38.8	Xtriage
Anisotropy	0.470	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 30.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 30831 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	7644	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.15% 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, 0.375 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: CA, 3AT

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	0.45	0/3876	0.61	0/5258
1	B	0.43	0/3837	0.62	2/5206 (0.0%)
All	All	0.44	0/7713	0.62	2/10464 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	415	VAL	CB-CA-C	-5.37	101.19	111.40
1	B	476	LEU	CA-CB-CG	5.13	127.11	115.30

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	3784	0	3791	52	0
1	B	3745	0	3745	78	0
2	A	30	0	11	1	0
2	B	30	0	11	5	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	27	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	26	0	0	5	0
All	All	7644	0	7558	130	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 (130) 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:461:SER:O	1:A:465:THR:OG1	2.07	0.71
1:B:180:ASN:ND2	4:B:713:HOH:O	2.17	0.71
1:B:177:ILE:HG23	1:B:181:LEU:HD12	1.74	0.69
1:B:74:TRP:O	1:B:77:ASP:N	2.25	0.68
1:A:94:GLY:H	1:A:177:ILE:HB	1.58	0.68
1:A:497:PRO:O	1:A:499:GLU:N	2.24	0.68
1:B:145:LEU:HD23	1:B:158:PHE:HB3	1.77	0.66
2:B:601:3AT:O2'	4:B:724:HOH:O	2.15	0.65
1:A:120:PRO:HG3	1:A:177:ILE:HG12	1.80	0.64
1:A:233:ARG:NH1	1:A:339:GLU:OE1	2.31	0.63
1:A:74:TRP:CE2	1:A:132:PHE:HB2	2.34	0.62
1:B:79:SER:HB3	1:B:89:VAL:HG13	1.82	0.62
1:B:487:VAL:HG13	1:B:491:GLN:HB2	1.82	0.62
1:A:376:LEU:HD21	1:A:462:PHE:HE2	1.65	0.61
1:B:220:ARG:O	1:B:224:ARG:HG3	2.01	0.61
1:B:232:ARG:NH1	1:B:361:LEU:O	2.36	0.58
1:B:200:LEU:O	1:B:204:ARG:HG3	2.04	0.58
1:B:53:ASP:HB2	1:B:56:GLU:H	1.69	0.57
1:B:101:SER:OG	2:B:601:3AT:O3G	2.10	0.57
1:A:103:ARG:HG2	1:A:206:THR:HG21	1.88	0.56
1:B:71:VAL:HG21	1:B:96:ILE:HD11	1.88	0.55
1:A:262:ALA:HB1	1:A:266:THR:HB	1.88	0.55
1:B:459:ILE:O	1:B:463:THR:HG23	2.06	0.55
1:A:121:ARG:HG3	1:A:174:ILE:O	2.07	0.55
1:B:189:LEU:HD22	1:B:197:ILE:HG12	1.88	0.54
1:B:120:PRO:HG3	1:B:177:ILE:HG12	1.90	0.54
1:B:374:ILE:HD11	1:B:455:LEU:HD13	1.90	0.54
1:A:129:PHE:CE2	1:A:149:GLU:HB3	2.43	0.54
1:B:207:ASP:O	1:B:211:HIS:ND1	2.40	0.54
1:A:232:ARG:NH1	1:A:361:LEU:O	2.41	0.53
1:B:94:GLY:H	1:B:177:ILE:HB	1.72	0.53
1:A:21:ILE:HG13	1:A:22:THR:HG23	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:462:PHE:O	1:B:466:VAL:HG23	2.09	0.53
1:B:92:VAL:HG13	1:B:120:PRO:HG2	1.90	0.53
1:A:205:VAL:HG21	2:A:601:3AT:H2	1.90	0.53
1:B:50:VAL:HG21	1:B:228:LEU:HG	1.89	0.53
1:A:103:ARG:HG2	1:A:206:THR:CG2	2.39	0.53
1:A:207:ASP:O	1:A:211:HIS:ND1	2.43	0.52
1:B:360:LEU:HB3	1:B:361:LEU:HD13	1.92	0.52
1:B:91:THR:HG23	1:B:176:THR:HG21	1.91	0.52
1:B:208:GLU:HG3	1:B:318:ALA:HB2	1.90	0.52
1:A:181:LEU:HD22	1:A:182:ASP:H	1.75	0.52
1:B:229:TRP:CD1	1:B:360:LEU:HD22	2.45	0.51
1:B:486:HIS:HD2	1:B:487:VAL:N	2.09	0.51
1:B:186:ASP:OD1	1:B:204:ARG:NH2	2.43	0.51
1:B:178:SER:OG	1:B:179:ASP:N	2.44	0.51
1:A:186:ASP:OD1	1:A:204:ARG:NH2	2.38	0.50
1:B:134:GLU:O	1:B:138:HIS:ND1	2.41	0.50
1:A:462:PHE:O	1:A:466:VAL:HG23	2.12	0.50
1:B:205:VAL:HG21	2:B:601:3AT:H2	1.92	0.50
1:B:476:LEU:HA	1:B:480:MET:CE	2.42	0.50
1:B:400:LEU:O	1:B:400:LEU:HD22	2.11	0.50
1:B:397:ILE:HG21	1:B:415:VAL:HG11	1.94	0.49
1:B:25:ILE:HD12	1:B:258:LEU:HB3	1.95	0.49
2:B:601:3AT:O2A	4:B:726:HOH:O	2.20	0.49
1:A:370:TYR:CD2	1:A:439:GLY:HA3	2.48	0.49
1:B:370:TYR:CD2	1:B:439:GLY:HA3	2.47	0.49
1:A:205:VAL:HG22	1:A:316:THR:HG23	1.95	0.49
1:B:350:ILE:HD13	1:B:357:TRP:HA	1.96	0.48
1:A:496:LEU:HA	1:A:497:PRO:HD3	1.73	0.48
1:A:45:MET:HG2	1:A:228:LEU:HD12	1.93	0.48
1:A:396:LYS:HE3	1:A:469:GLN:OE1	2.14	0.48
1:A:477:LYS:H	1:A:480:MET:HE2	1.78	0.48
1:B:130:GLN:HG3	1:B:131:SER:H	1.79	0.48
1:A:229:TRP:O	1:A:233:ARG:HB2	2.14	0.48
1:B:121:ARG:HG3	1:B:122:HIS:CD2	2.50	0.47
1:B:476:LEU:HA	1:B:480:MET:HE1	1.96	0.47
1:B:205:VAL:O	1:B:209:ILE:HG13	2.15	0.47
1:B:435:MET:HE1	1:B:495:TYR:HD2	1.80	0.47
1:B:57:LEU:O	1:B:61:LEU:HD23	2.15	0.46
1:A:185:ASP:OD2	1:A:187:SER:OG	2.21	0.46
1:A:79:SER:HB2	1:A:122:HIS:CE1	2.51	0.46
1:A:204:ARG:NH1	1:A:319:TYR:O	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:342:GLN:O	1:A:346:VAL:HG23	2.17	0.45
1:A:371:ARG:HA	1:A:371:ARG:HD2	1.67	0.45
1:B:234:GLY:O	4:B:725:HOH:O	2.21	0.45
1:A:243:LEU:N	1:A:336:MET:HE1	2.32	0.45
1:A:499:GLU:O	1:A:501:LEU:N	2.49	0.45
1:B:232:ARG:HA	1:B:232:ARG:HD2	1.57	0.45
1:B:106:VAL:HG21	1:B:223:LEU:CD1	2.47	0.45
1:A:145:LEU:HD23	1:A:145:LEU:HA	1.79	0.44
1:B:123:VAL:O	1:B:171:ARG:HG3	2.17	0.44
1:A:258:LEU:HD23	1:A:258:LEU:HA	1.79	0.44
1:B:105:GLY:HA2	4:B:717:HOH:O	2.16	0.44
1:B:243:LEU:HD13	1:B:336:MET:HE1	1.99	0.44
1:B:288:GLN:HA	1:B:289:PRO:HD3	1.77	0.44
1:A:475:MET:O	1:A:477:LYS:HG3	2.18	0.44
1:A:29:SER:O	1:B:406:ARG:HG2	2.17	0.44
1:B:92:VAL:HA	1:B:176:THR:HG22	2.00	0.44
1:A:74:TRP:CE3	1:A:75:ILE:HG13	2.53	0.44
1:B:181:LEU:HD22	1:B:183:LEU:HD23	2.00	0.44
1:A:496:LEU:HA	1:A:496:LEU:HD12	1.71	0.44
1:A:121:ARG:NH1	1:A:175:GLN:OE1	2.48	0.44
1:B:30:PRO:HD3	1:B:261:ASN:O	2.18	0.44
1:B:120:PRO:HB3	1:B:175:GLN:O	2.18	0.43
1:B:496:LEU:HD12	1:B:496:LEU:HA	1.85	0.43
1:B:235:ILE:HB	1:B:336:MET:HE3	2.00	0.43
1:B:19:TYR:CE2	1:B:214:PRO:HA	2.53	0.43
1:B:92:VAL:HG11	1:B:122:HIS:HE1	1.83	0.43
1:A:128:PHE:CE2	1:A:156:ILE:HG13	2.54	0.43
1:B:181:LEU:HD21	1:B:188:ARG:HH11	1.83	0.43
1:B:125:ARG:O	1:B:129:PHE:HD2	2.02	0.42
1:A:185:ASP:O	1:A:188:ARG:HG3	2.19	0.42
1:A:61:LEU:HD13	4:A:704:HOH:O	2.20	0.42
1:B:246:VAL:HG22	2:B:601:3AT:H2'1	2.02	0.42
1:A:145:LEU:HD23	1:A:158:PHE:HB3	2.01	0.42
1:A:67:LEU:HD13	1:A:96:ILE:CD1	2.49	0.42
1:B:233:ARG:HA	1:B:364:PRO:HD3	2.01	0.42
1:B:268:VAL:O	1:B:271:PHE:HB3	2.19	0.42
1:B:396:LYS:O	1:B:399:VAL:HB	2.20	0.42
1:B:259:TYR:CZ	1:B:270:LYS:HD2	2.54	0.42
1:B:486:HIS:CD2	1:B:487:VAL:N	2.87	0.41
1:A:103:ARG:O	1:A:103:ARG:HG3	2.18	0.41
1:B:366:PHE:O	1:B:373:TYR:OH	2.33	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:459:ILE:HD13	1:B:484:ALA:O	2.20	0.41
1:B:316:THR:HA	1:B:317:PRO:HD3	1.96	0.41
1:B:154:PRO:O	1:B:168:VAL:HG22	2.19	0.41
1:A:193:ASP:O	1:A:196:CYS:N	2.54	0.41
1:B:243:LEU:N	1:B:336:MET:HE1	2.35	0.41
1:B:99:PHE:CZ	1:B:114:ASP:HB3	2.56	0.41
1:B:382:THR:OG1	1:B:383:GLU:N	2.53	0.41
1:A:487:VAL:HG12	1:A:488:LYS:O	2.21	0.41
1:A:335:VAL:HG21	1:A:414:HIS:CD2	2.56	0.41
1:A:74:TRP:CG	1:A:132:PHE:HD1	2.39	0.40
1:B:184:ARG:HD3	1:B:208:GLU:OE2	2.21	0.40
1:A:28:ALA:HB1	1:B:406:ARG:HB3	2.03	0.40
1:B:213:VAL:HA	1:B:214:PRO:HD2	1.96	0.40
1:B:112:ASP:HB3	1:B:164:GLU:HB3	2.04	0.40
1:B:133:PHE:CZ	1:B:145:LEU:HD13	2.56	0.40
1:A:312:MET:N	1:A:313:PRO:HD3	2.37	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	
1	A	461/514 (90%)	425 (92%)	31 (7%)	5 (1%)	17	50
1	B	456/514 (89%)	427 (94%)	23 (5%)	6 (1%)	15	44
All	All	917/1028 (89%)	852 (93%)	54 (6%)	11 (1%)	16	47

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	500	ILE
1	B	496	LEU

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Mol	Chain	Res	Type
1	B	237	SER
1	B	490	LYS
1	A	237	SER
1	A	497	PRO
1	A	498	ALA
1	B	75	ILE
1	B	421	PRO
1	A	303	VAL
1	B	148	VAL

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	420/465 (90%)	368 (88%)	52 (12%)	6	17
1	B	416/465 (90%)	371 (89%)	45 (11%)	8	23
All	All	836/930 (90%)	739 (88%)	97 (12%)	7	20

All (97) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	41	LEU
1	A	61	LEU
1	A	67	LEU
1	A	73	GLU
1	A	96	ILE
1	A	103	ARG
1	A	130	GLN
1	A	144	ASN
1	A	145	LEU
1	A	146	ARG
1	A	160	PHE
1	A	168	VAL
1	A	181	LEU
1	A	184	ARG

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Mol	Chain	Res	Type
1	A	207	ASP
1	A	213	VAL
1	A	223	LEU
1	A	228	LEU
1	A	232	ARG
1	A	233	ARG
1	A	239	MET
1	A	254	ARG
1	A	273	LEU
1	A	288	GLN
1	A	306	SER
1	A	316	THR
1	A	321	GLN
1	A	330	THR
1	A	342	GLN
1	A	360	LEU
1	A	361	LEU
1	A	369	LYS
1	A	371	ARG
1	A	376	LEU
1	A	379	SER
1	A	381	SER
1	A	383	GLU
1	A	384	GLU
1	A	400	LEU
1	A	411	THR
1	A	433	VAL
1	A	435	MET
1	A	443	ARG
1	A	455	LEU
1	A	461	SER
1	A	468	ARG
1	A	473	ILE
1	A	478	GLU
1	A	491	GLN
1	A	492	LEU
1	A	499	GLU
1	A	500	ILE
1	B	29	SER
1	B	41	LEU
1	B	67	LEU
1	B	68	ASN

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Mol	Chain	Res	Type
1	B	69	ASN
1	B	79	SER
1	B	84	LEU
1	B	103	ARG
1	B	121	ARG
1	B	124	GLU
1	B	145	LEU
1	B	146	ARG
1	B	160	PHE
1	B	181	LEU
1	B	182	ASP
1	B	184	ARG
1	B	217	GLU
1	B	223	LEU
1	B	228	LEU
1	B	232	ARG
1	B	239	MET
1	B	246	VAL
1	B	254	ARG
1	B	273	LEU
1	B	290	GLU
1	B	316	THR
1	B	321	GLN
1	B	342	GLN
1	B	355	SER
1	B	360	LEU
1	B	361	LEU
1	B	376	LEU
1	B	382	THR
1	B	383	GLU
1	B	384	GLU
1	B	387	LEU
1	B	390	VAL
1	B	392	LEU
1	B	400	LEU
1	B	415	VAL
1	B	418	GLN
1	B	456	THR
1	B	478	GLU
1	B	490	LYS
1	B	496	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	493	HIS
1	A	494	HIS
1	B	122	HIS
1	B	486	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](#)

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	3AT	A	601	3	24,32,32	1.50	2 (8%)	26,50,50	2.42	6 (23%)
2	3AT	B	601	3	24,32,32	1.60	4 (16%)	26,50,50	2.55	6 (23%)

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	3AT	A	601	3	-	0/18/34/34	0/3/3/3
2	3AT	B	601	3	-	0/18/34/34	0/3/3/3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	601	3AT	O2'-C2'	-4.55	1.33	1.43
2	A	601	3AT	O2'-C2'	-4.16	1.34	1.43
2	B	601	3AT	C2'-C1'	-2.02	1.49	1.53
2	B	601	3AT	O4'-C4'	2.17	1.49	1.44
2	A	601	3AT	C6-N6	2.60	1.43	1.34
2	B	601	3AT	C6-N6	2.82	1.43	1.34

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	3AT	N3-C2-N1	-10.56	120.81	128.89
2	A	601	3AT	N3-C2-N1	-10.05	121.20	128.89
2	B	601	3AT	PB-O3B-PG	-2.90	122.93	132.67
2	B	601	3AT	PA-O3A-PB	-2.62	125.36	132.73
2	B	601	3AT	C4-C5-N7	-2.50	107.18	109.48
2	A	601	3AT	PA-O3A-PB	-2.27	126.36	132.73
2	A	601	3AT	C2-N1-C6	2.08	122.48	118.77
2	A	601	3AT	O4'-C4'-C5'	2.19	112.77	109.54
2	B	601	3AT	O4'-C4'-C5'	2.22	112.82	109.54
2	A	601	3AT	O3A-PA-O5'	2.73	110.17	102.94
2	A	601	3AT	O4'-C1'-N9	2.98	114.33	108.10
2	B	601	3AT	O4'-C1'-N9	3.50	115.42	108.10

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	3AT	1	0
2	B	601	3AT	5	0

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	467/514 (90%)	-0.28	5 (1%) 82 74	5, 21, 62, 99	2 (0%)
1	B	462/514 (89%)	-0.24	4 (0%) 85 79	6, 23, 59, 92	2 (0%)
All	All	929/1028 (90%)	-0.26	9 (0%) 84 77	5, 22, 62, 99	4 (0%)

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	152	PHE	3.0
1	B	497	PRO	3.0
1	A	86	PRO	2.9
1	A	457	TYR	2.6
1	A	431	ASN	2.5
1	A	498	ALA	2.3
1	B	191	SER	2.1
1	A	82	LYS	2.1
1	B	87	SER	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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	3AT	A	601	30/30	0.96	0.17	2.00	11,24,29,37	7
2	3AT	B	601	30/30	0.96	0.15	0.60	16,25,37,48	7
3	CA	B	602	1/1	0.98	0.05	-	34,34,34,34	0
3	CA	A	602	1/1	0.95	0.07	-	31,31,31,31	0

6.5 Other polymers

There are no such residues in this entry.