Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 09:04 AM GMT

PDB ID	:	3H4B
Title	:	Ternary complex of human DNA polymerase iota with template U/T and incoming dATP
Authors	:	Jain, R.; Nair, D.T.; Johnson, R.E.; Prakash, L.; Prakash, S.; Aggarwal, A.K.
Deposited on	:	2009-04-18
Resolution	:	2.85 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:

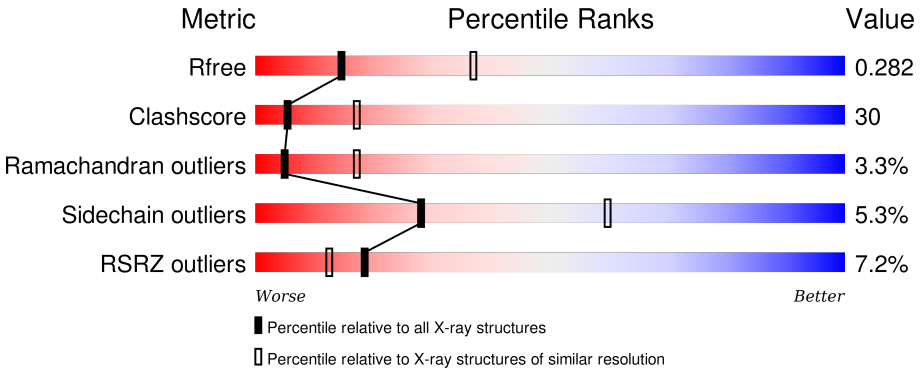
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 2.85 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	2228 (2.90-2.82)
Clashscore	102246	2499 (2.90-2.82)
Ramachandran outliers	100387	2439 (2.90-2.82)
Sidechain outliers	100360	2442 (2.90-2.82)
RSRZ outliers	91569	2236 (2.90-2.82)

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	A	390	7%, 52%, 38%, . . .
2	P	7	14%, 71%, 14%
3	T	9	78%, 22%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 3404 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 iota.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	373	Total	C	N	O	S	0	0	0
			2877	1809	504	543	21			

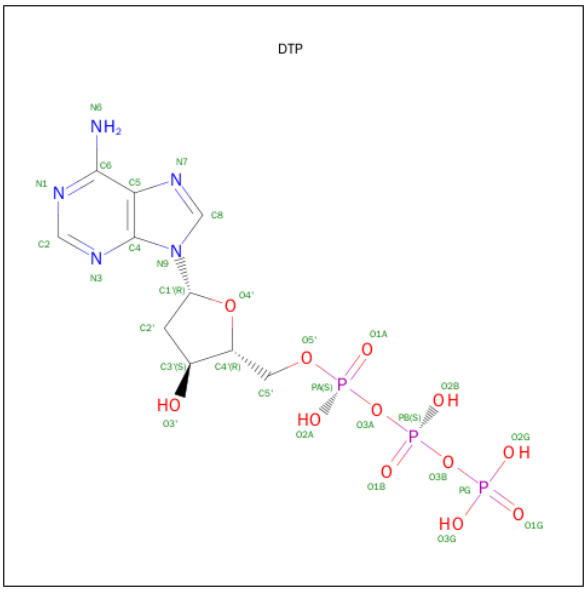
- Molecule 2 is a DNA chain called 5'-D(*AP*GP*GP*AP*CP*CP*(DOC))-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	P	7	Total	C	N	O	P	0	0	0
			139	67	29	37	6			

- Molecule 3 is a DNA chain called 5'-D(*TP*(BRU)P*GP*GP*GP*TP*CP*CP*T)-3'.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	T	9	Total	Br	C	N	O	P	0	1	0
			201	2	96	31	63	9			

- Molecule 4 is 2'-DEOXYADENOSINE 5'-TRIPHOSPHATE (three-letter code: DTP) (formula: $C_{10}H_{16}N_5O_{12}P_3$).



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

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

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

- Molecule 6 is water.

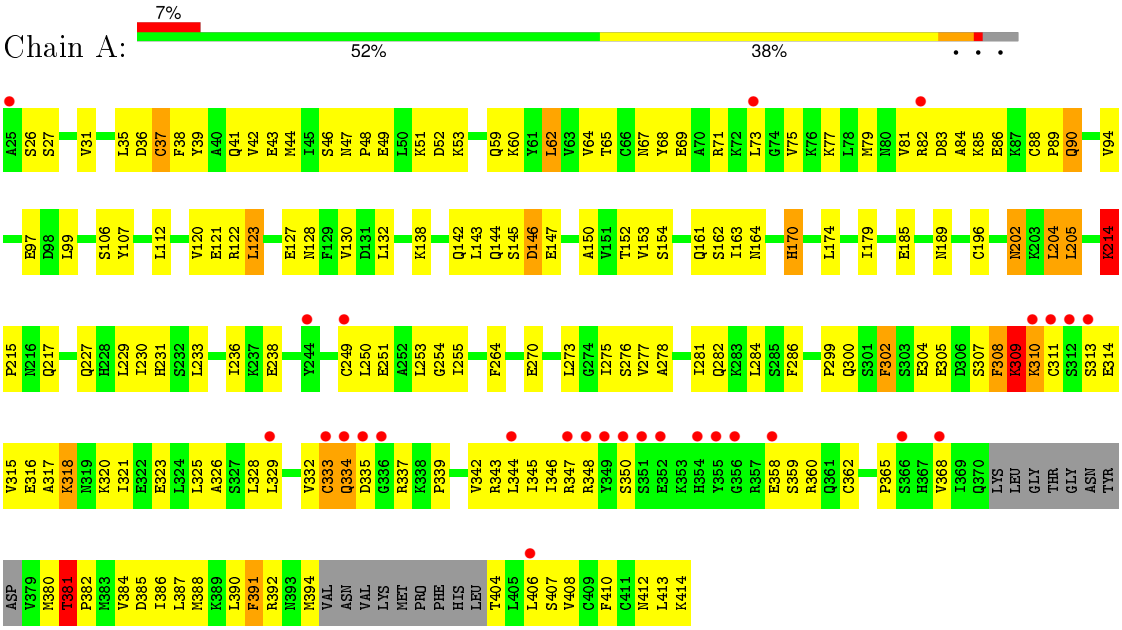
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	141	Total	O	0	0
			141	141		
6	P	7	Total	O	0	0
			7	7		
6	T	8	Total	O	0	0
			8	8		

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.

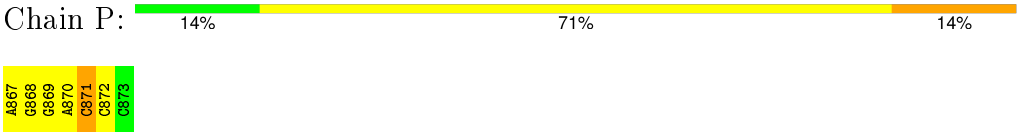
- Molecule 1: DNA polymerase iota

Chain A:



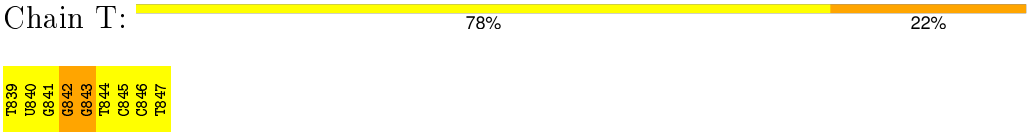
- Molecule 2: 5'-D(*AP*GP*GP*AP*CP*CP*(DOC))-3'

Chain P:



- Molecule 3: 5'-D(*TP*(BRU)P*GP*GP*GP*TP*CP*CP*T)-3'

Chain T:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α, β, γ	98.00Å 98.00Å 203.67Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	39.73 – 2.85 41.54 – 2.85	Depositor EDS
% Data completeness (in resolution range)	96.5 (39.73-2.85) 96.6 (41.54-2.85)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	5.91 (at 2.86Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.240 , 0.281 0.240 , 0.282	Depositor DCC
R_{free} test set	1094 reflections (8.69%)	DCC
Wilson B-factor (Å^2)	45.5	Xtriage
Anisotropy	0.165	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 63.3	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 14175 reflections	Xtriage
F_o,F_c correlation	0.90	EDS
Total number of atoms	3404	wwPDB-VP
Average B, all atoms (Å^2)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.51% 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, MG, DTP, BRU

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	0.37	0/2915	0.64	1/3934 (0.0%)
2	P	0.47	0/136	0.77	0/208
3	T	0.50	0/178	0.90	0/271
All	All	0.38	0/3229	0.66	1/4413 (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	P	0	2
3	T	0	2
All	All	0	4

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	214	LYS	N-CA-C	6.24	127.86	111.00

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	P	871	DC	Sidechain
2	P	872	DC	Sidechain
3	T	842	DG	Sidechain
3	T	843	DG	Sidechain

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	2877	0	2899	163	0
2	P	139	0	79	4	0
3	T	201	0	111	29	0
4	A	30	0	10	3	0
5	A	1	0	0	0	0
6	A	141	0	0	9	0
6	P	7	0	0	1	0
6	T	8	0	0	2	0
All	All	3404	0	3099	189	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 30.

All (189) 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:44:MET:HE2	1:A:51:LYS:HA	1.41	1.01
3:T:842:DG:H2"	3:T:843:DG:H5"	1.47	0.95
1:A:77:LYS:HB2	1:A:77:LYS:NZ	1.87	0.89
1:A:344:LEU:HD11	1:A:387:LEU:HD22	1.53	0.87
3:T:842:DG:H2"	3:T:843:DG:C5'	2.06	0.85
1:A:309:LYS:H	1:A:309:LYS:HD3	1.39	0.84
1:A:365:PRO:O	1:A:368:VAL:HG22	1.77	0.84
1:A:313:SER:O	1:A:314:GLU:HB2	1.78	0.81
1:A:388:MET:O	1:A:391:PHE:HB3	1.80	0.81
3:T:839:DT:H2"	3:T:840[A]:BRU:O5'	1.82	0.79
1:A:270:GLU:HG3	1:A:275:ILE:HA	1.65	0.78
1:A:59:GLN:OE1	3:T:840[B]:BRU:H6	1.86	0.75
1:A:75:VAL:HA	1:A:79:MET:HE2	1.68	0.75
1:A:321:ILE:O	1:A:325:LEU:HD13	1.87	0.75
1:A:44:MET:CE	1:A:67:ASN:HD22	1.99	0.75
1:A:347:ARG:HD2	1:A:404:THR:OG1	1.87	0.75
1:A:309:LYS:H	1:A:309:LYS:CD	1.96	0.74
1:A:164:ASN:H	1:A:170:HIS:HD2	1.36	0.73
1:A:304:GLU:HG3	1:A:328:LEU:HG	1.69	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:T:841:DG:H2"	3:T:842:DG:H5"	1.69	0.72
1:A:384:VAL:O	1:A:388:MET:HG2	1.89	0.71
1:A:77:LYS:HB2	1:A:77:LYS:HZ2	1.54	0.70
1:A:344:LEU:HD21	1:A:387:LEU:HB3	1.71	0.70
1:A:39:TYR:HB2	1:A:65:THR:HG21	1.74	0.69
3:T:839:DT:H2'	3:T:840[A]:BRU:BR	2.48	0.68
1:A:254:GLY:O	6:A:469:HOH:O	2.12	0.68
1:A:318:LYS:HD2	1:A:388:MET:HE1	1.76	0.67
3:T:841:DG:C2'	3:T:842:DG:H5"	2.24	0.67
1:A:51:LYS:O	1:A:53:LYS:N	2.28	0.67
1:A:77:LYS:HE2	4:A:875:DTP:PB	2.34	0.67
4:A:875:DTP:O1G	6:A:526:HOH:O	2.12	0.67
1:A:310:LYS:HG3	6:A:507:HOH:O	1.94	0.67
1:A:79:MET:CE	1:A:84:ALA:HA	2.25	0.67
1:A:44:MET:HE1	1:A:67:ASN:HD22	1.59	0.66
1:A:365:PRO:HB2	1:A:368:VAL:HG13	1.78	0.66
1:A:335:ASP:OD2	1:A:337:ARG:HD3	1.96	0.66
3:T:841:DG:H1'	3:T:842:DG:H5"	1.78	0.65
1:A:71:ARG:HH22	4:A:875:DTP:PG	2.19	0.65
1:A:233:LEU:HD22	1:A:238:GLU:HB3	1.79	0.65
1:A:305:GLU:HG3	1:A:407:SER:HB3	1.78	0.65
1:A:132:LEU:HD22	1:A:179:ILE:HG21	1.79	0.64
2:P:870:DA:H1'	2:P:871:DC:H5'	1.79	0.64
1:A:325:LEU:HD11	1:A:387:LEU:CD1	2.27	0.64
1:A:320:LYS:O	1:A:323:GLU:HB2	1.97	0.64
2:P:867:DA:H1'	2:P:868:DG:H5'	1.80	0.64
1:A:315:VAL:C	1:A:317:ALA:H	2.00	0.63
1:A:381:THR:O	6:A:517:HOH:O	2.16	0.63
1:A:69:GLU:O	1:A:73:LEU:HD13	1.99	0.63
1:A:387:LEU:HA	1:A:390:LEU:HD12	1.81	0.62
1:A:381:THR:OG1	1:A:382:PRO:HD3	1.98	0.62
1:A:325:LEU:O	1:A:329:LEU:HD13	1.99	0.62
1:A:62:LEU:CD1	3:T:840[A]:BRU:H5"	2.30	0.62
1:A:88:CYS:SG	1:A:90:GLN:HG3	2.40	0.62
1:A:143:LEU:HA	1:A:147:GLU:OE2	1.98	0.62
1:A:43:GLU:HG3	1:A:94:VAL:HG21	1.81	0.61
1:A:214:LYS:HB3	1:A:215:PRO:HD3	1.83	0.61
1:A:51:LYS:C	1:A:53:LYS:H	2.04	0.60
1:A:164:ASN:H	1:A:170:HIS:CD2	2.18	0.60
1:A:202:ASN:ND2	1:A:205:LEU:H	2.00	0.59
2:P:870:DA:N3	6:P:80:HOH:O	2.32	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:300:GLN:NE2	6:A:491:HOH:O	2.35	0.58
1:A:196:CYS:SG	1:A:214:LYS:O	2.62	0.57
1:A:249:CYS:SG	1:A:273:LEU:HD21	2.44	0.57
1:A:107:TYR:OH	1:A:299:PRO:HG3	2.05	0.57
3:T:839:DT:H5"	6:T:170:HOH:O	2.04	0.57
1:A:62:LEU:HD11	3:T:839:DT:O2	2.05	0.57
1:A:44:MET:HE3	1:A:67:ASN:HD22	1.69	0.56
1:A:59:GLN:OE1	1:A:64:VAL:HG11	2.05	0.56
1:A:81:VAL:HG12	1:A:85:LYS:HE3	1.87	0.56
1:A:85:LYS:O	1:A:89:PRO:HG3	2.06	0.56
1:A:227:GLN:O	1:A:230:ILE:HG22	2.05	0.56
1:A:325:LEU:HD11	1:A:387:LEU:HD11	1.88	0.56
1:A:79:MET:HE3	1:A:84:ALA:HA	1.87	0.56
3:T:839:DT:H2"	3:T:840[B]:BRU:O5'	2.06	0.55
1:A:35:LEU:HD13	6:A:436:HOH:O	2.05	0.55
1:A:79:MET:HE1	1:A:84:ALA:HA	1.89	0.54
1:A:332:VAL:CG1	1:A:339:PRO:HD3	2.37	0.54
1:A:27:SER:HA	6:A:498:HOH:O	2.08	0.54
1:A:62:LEU:HD11	3:T:839:DT:C2	2.42	0.54
1:A:62:LEU:HD12	3:T:840[A]:BRU:H5"	1.89	0.54
1:A:385:ASP:HA	1:A:388:MET:HG2	1.89	0.54
1:A:163:ILE:HD13	1:A:174:LEU:HD11	1.90	0.54
1:A:339:PRO:HG3	1:A:410:PHE:HD2	1.74	0.53
1:A:250:LEU:HA	1:A:253:LEU:HD12	1.89	0.53
2:P:868:DG:H2"	2:P:869:DG:OP2	2.07	0.53
1:A:343:ARG:C	1:A:344:LEU:HD12	2.28	0.53
1:A:82:ARG:HG2	6:A:490:HOH:O	2.08	0.53
1:A:325:LEU:CD1	1:A:408:VAL:HG21	2.38	0.53
1:A:388:MET:HE3	1:A:388:MET:HA	1.89	0.53
1:A:204:LEU:HG	1:A:284:LEU:HD22	1.89	0.53
1:A:215:PRO:O	1:A:217:GLN:HG3	2.09	0.52
1:A:365:PRO:HB2	1:A:368:VAL:CG1	2.40	0.52
1:A:106:SER:OG	1:A:122:ARG:NH2	2.43	0.52
1:A:318:LYS:CD	1:A:388:MET:HE1	2.39	0.52
1:A:79:MET:HE3	1:A:84:ALA:CB	2.40	0.52
1:A:386:ILE:O	1:A:390:LEU:HG	2.10	0.51
1:A:344:LEU:O	1:A:359:SER:HA	2.11	0.51
3:T:842:DG:C2'	3:T:843:DG:H5"	2.31	0.51
1:A:270:GLU:HA	1:A:278:ALA:CB	2.41	0.51
3:T:845:DC:H2"	3:T:846:DC:OP2	2.11	0.50
1:A:161:GLN:HG2	1:A:162:SER:N	2.27	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:99:LEU:HD12	3:T:842:DG:H4'	1.94	0.50
1:A:347:ARG:HG2	1:A:348:ARG:N	2.27	0.50
1:A:185:GLU:OE2	1:A:189:ASN:ND2	2.45	0.49
1:A:309:LYS:O	1:A:310:LYS:C	2.51	0.49
1:A:343:ARG:HG2	1:A:344:LEU:N	2.26	0.49
3:T:841:DG:C1'	3:T:842:DG:H5"	2.41	0.49
1:A:309:LYS:N	1:A:309:LYS:HD3	2.18	0.49
1:A:144:GLN:O	1:A:146:ASP:N	2.37	0.49
1:A:36:ASP:O	1:A:37:CYS:C	2.50	0.49
1:A:360:ARG:HG2	1:A:394:MET:SD	2.52	0.49
1:A:144:GLN:H	1:A:147:GLU:CD	2.16	0.49
3:T:841:DG:H2"	3:T:842:DG:C5'	2.42	0.49
1:A:385:ASP:HA	1:A:388:MET:CG	2.43	0.48
1:A:82:ARG:HG3	1:A:83:ASP:H	1.77	0.48
1:A:112:LEU:HD23	1:A:112:LEU:C	2.34	0.48
1:A:153:VAL:HG22	1:A:154:SER:N	2.28	0.48
1:A:123:LEU:O	1:A:127:GLU:HB2	2.13	0.48
1:A:230:ILE:CG2	1:A:231:HIS:N	2.77	0.48
1:A:138:LYS:O	1:A:142:GLN:HG2	2.14	0.48
1:A:202:ASN:OD1	1:A:205:LEU:HD22	2.14	0.48
1:A:302:PHE:N	1:A:302:PHE:CD1	2.82	0.48
1:A:308:PHE:O	1:A:310:LYS:N	2.47	0.47
1:A:233:LEU:HD22	1:A:238:GLU:CB	2.43	0.47
1:A:85:LYS:HB2	1:A:85:LYS:NZ	2.29	0.47
1:A:79:MET:HE3	1:A:84:ALA:CA	2.45	0.47
1:A:46:SER:O	1:A:48:PRO:HD3	2.15	0.47
1:A:316:GLU:C	1:A:318:LYS:H	2.17	0.46
1:A:122:ARG:NH1	6:A:13:HOH:O	2.48	0.46
1:A:153:VAL:CB	1:A:174:LEU:HD22	2.46	0.46
1:A:236:ILE:HD12	1:A:250:LEU:HD12	1.97	0.46
1:A:144:GLN:H	1:A:147:GLU:CG	2.28	0.46
1:A:214:LYS:HB3	1:A:215:PRO:CD	2.46	0.46
1:A:412:ASN:OD1	1:A:414:LYS:HE3	2.16	0.46
1:A:31:VAL:CG1	1:A:130:VAL:HB	2.45	0.46
1:A:323:GLU:O	1:A:326:ALA:HB3	2.16	0.46
1:A:75:VAL:HA	1:A:79:MET:CE	2.42	0.45
1:A:62:LEU:CD1	3:T:840[B]:BRU:H5"	2.45	0.45
1:A:214:LYS:CB	1:A:215:PRO:HD3	2.44	0.45
1:A:47:ASN:OD1	1:A:49:GLU:HB2	2.16	0.45
1:A:315:VAL:C	1:A:317:ALA:N	2.68	0.45
3:T:846:DC:H1'	3:T:847:DT:H5"	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:39:TYR:HB2	1:A:65:THR:CG2	2.43	0.45
1:A:202:ASN:HD21	1:A:205:LEU:H	1.64	0.45
1:A:120:VAL:HG22	1:A:130:VAL:HG22	2.00	0.44
1:A:325:LEU:HD11	1:A:408:VAL:HG21	1.99	0.44
1:A:236:ILE:HD12	1:A:250:LEU:CD1	2.48	0.44
1:A:346:ILE:HG22	1:A:406:LEU:CD2	2.47	0.44
1:A:337:ARG:HB3	1:A:414:LYS:O	2.18	0.44
1:A:277:VAL:O	1:A:281:ILE:HG12	2.17	0.44
1:A:97:GLU:HG2	3:T:841:DG:OP1	2.17	0.44
1:A:345:ILE:HA	1:A:358:GLU:O	2.17	0.44
1:A:79:MET:HE3	1:A:84:ALA:HB2	2.00	0.44
3:T:847:DT:H6	3:T:847:DT:H5'	1.83	0.44
3:T:841:DG:N2	6:T:135:HOH:O	2.51	0.43
1:A:304:GLU:O	1:A:407:SER:HB2	2.18	0.43
1:A:282:GLN:O	1:A:286:PHE:HD1	2.01	0.43
1:A:381:THR:O	1:A:382:PRO:C	2.57	0.43
1:A:38:PHE:O	1:A:42:VAL:HG23	2.19	0.43
1:A:309:LYS:O	1:A:311:CYS:N	2.51	0.43
1:A:380:MET:O	1:A:381:THR:C	2.57	0.43
1:A:413:LEU:N	1:A:413:LEU:HD22	2.33	0.43
3:T:843:DG:H2'	3:T:844:DT:H72	2.01	0.43
3:T:846:DC:H2"	3:T:847:DT:H5'	2.01	0.43
1:A:153:VAL:HB	1:A:174:LEU:HD22	2.01	0.42
1:A:202:ASN:HD22	1:A:202:ASN:C	2.23	0.42
1:A:333:CYS:O	1:A:334:GLN:C	2.57	0.42
1:A:86:GLU:O	1:A:86:GLU:HG3	2.18	0.42
1:A:60:LYS:HE2	1:A:307:SER:O	2.18	0.42
1:A:82:ARG:C	1:A:84:ALA:N	2.73	0.42
1:A:313:SER:OG	1:A:315:VAL:O	2.38	0.42
3:T:842:DG:H2"	3:T:843:DG:H5'	1.94	0.42
1:A:315:VAL:O	1:A:317:ALA:N	2.47	0.42
1:A:68:TYR:OH	1:A:214:LYS:HD2	2.20	0.42
1:A:36:ASP:O	1:A:38:PHE:N	2.53	0.42
1:A:121:GLU:O	1:A:128:ASN:HA	2.20	0.42
3:T:843:DG:C8	3:T:844:DT:H72	2.55	0.42
1:A:196:CYS:HA	1:A:217:GLN:O	2.20	0.42
3:T:839:DT:C2'	3:T:840[A]:BRU:O5'	2.61	0.41
1:A:251:GLU:C	1:A:253:LEU:N	2.73	0.41
1:A:388:MET:CE	1:A:391:PHE:HD1	2.33	0.41
1:A:255:ILE:HD11	1:A:264:PHE:CG	2.56	0.41
1:A:36:ASP:O	1:A:41:GLN:NE2	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:123:LEU:HD23	1:A:123:LEU:HA	1.73	0.41
1:A:150:ALA:O	1:A:152:THR:HG23	2.21	0.41
1:A:251:GLU:C	1:A:253:LEU:H	2.24	0.41
1:A:392:ARG:C	1:A:394:MET:N	2.75	0.41
1:A:413:LEU:N	1:A:413:LEU:CD2	2.85	0.40
1:A:380:MET:O	1:A:384:VAL:HG23	2.21	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
1	A	367/390 (94%)	307 (84%)	48 (13%)	12 (3%)	5 17

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	52	ASP
1	A	309	LYS
1	A	350	SER
1	A	26	SER
1	A	37	CYS
1	A	145	SER
1	A	310	LYS
1	A	334	GLN
1	A	146	ASP
1	A	276	SER
1	A	333	CYS
1	A	381	THR

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	321/354 (91%)	304 (95%)	17 (5%)	28 60

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	62	LEU
1	A	90	GLN
1	A	123	LEU
1	A	170	HIS
1	A	202	ASN
1	A	204	LEU
1	A	205	LEU
1	A	214	LYS
1	A	229	LEU
1	A	302	PHE
1	A	308	PHE
1	A	309	LYS
1	A	318	LYS
1	A	342	VAL
1	A	362	CYS
1	A	381	THR
1	A	391	PHE

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

Mol	Chain	Res	Type
1	A	58	GLN
1	A	67	ASN
1	A	90	GLN
1	A	170	HIS
1	A	202	ASN
1	A	235	HIS
1	A	256	ASN
1	A	262	GLN

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Mol	Chain	Res	Type
1	A	279	GLN
1	A	334	GLN
1	A	340	HIS

5.3.3 RNA

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains

3 non-standard protein/DNA/RNA residues are 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	873	3,2	11,19,20	0.80	0	14,26,29	1.13	2 (14%)
3	BRU	T	840[A]	3	13,21,22	1.58	3 (23%)	16,30,33	4.00	2 (12%)
3	BRU	T	840[B]	3	13,21,22	1.69	3 (23%)	16,30,33	4.14	2 (12%)

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	873	3,2	-	0/3/18/19	0/2/2/2
3	BRU	T	840[A]	3	-	0/3/21/22	0/2/2/2
3	BRU	T	840[B]	3	-	0/3/21/22	0/2/2/2

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	T	840[A]	BRU	C6-N1	2.31	1.38	1.35
3	T	840[B]	BRU	C6-N1	2.33	1.38	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	T	840[A]	BRU	C4-N3	2.70	1.38	1.33
3	T	840[B]	BRU	C4-N3	2.91	1.38	1.33
3	T	840[A]	BRU	C4-C5	3.95	1.43	1.38
3	T	840[B]	BRU	C4-C5	4.38	1.44	1.38

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	T	840[B]	BRU	C5-C4-N3	-8.69	114.72	124.00
3	T	840[A]	BRU	C5-C4-N3	-8.41	115.02	124.00
2	P	873	DOC	C2'-C1'-N1	2.10	116.73	112.49
2	P	873	DOC	C2-N3-C4	3.02	119.87	115.61
3	T	840[A]	BRU	C4-N3-C2	13.42	126.85	115.25
3	T	840[B]	BRU	C4-N3-C2	13.70	127.09	115.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	T	840[A]	BRU	5	0
3	T	840[B]	BRU	3	0

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
4	DTP	A	875	5	24,32,32	1.35	3 (12%)	32,50,50	1.41	8 (25%)

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	DTP	A	875	5	-	0/18/34/34	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	875	DTP	PA-O1A	-2.00	1.43	1.51
4	A	875	DTP	C2-N3	2.12	1.36	1.32
4	A	875	DTP	C3'-C4'	4.69	1.66	1.53

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	A	875	DTP	O4'-C4'-C3'	-2.70	98.88	105.67
4	A	875	DTP	C2'-C3'-C4'	-2.43	97.73	102.77
4	A	875	DTP	O4'-C4'-C5'	-2.29	101.12	109.32
4	A	875	DTP	PB-O3B-PG	-2.29	124.99	132.67
4	A	875	DTP	C5'-C4'-C3'	2.19	128.50	114.64
4	A	875	DTP	O3G-PG-O2G	2.28	116.06	107.38
4	A	875	DTP	O3A-PA-O5'	2.58	109.77	102.94
4	A	875	DTP	O2B-PB-O3A	2.76	117.62	105.09

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	875	DTP	3	0

5.7 Other polymers (i)

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

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	A	373/390 (95%)	0.25	28 (7%) 17 11	16, 48, 99, 120	1 (0%)
2	P	6/7 (85%)	0.30	0 100 100	38, 52, 54, 56	0
3	T	8/9 (88%)	0.07	0 100 100	30, 35, 42, 77	0
All	All	387/406 (95%)	0.24	28 (7%) 18 12	16, 48, 97, 120	1 (0%)

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	355	TYR	6.3
1	A	354	HIS	6.3
1	A	350	SER	4.6
1	A	348	ARG	4.5
1	A	311	CYS	4.3
1	A	25	ALA	4.1
1	A	313	SER	4.0
1	A	349	TYR	3.6
1	A	310	LYS	3.6
1	A	406	LEU	3.3
1	A	336	GLY	3.1
1	A	312	SER	3.0
1	A	352	GLU	2.9
1	A	366	SER	2.9
1	A	368	VAL	2.7
1	A	356	GLY	2.7
1	A	82	ARG	2.6
1	A	244	TYR	2.6
1	A	344	LEU	2.5
1	A	73	LEU	2.4
1	A	333	CYS	2.3
1	A	358	GLU	2.3
1	A	335	ASP	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	249	CYS	2.3
1	A	351	SER	2.2
1	A	334	GLN	2.2
1	A	347	ARG	2.2
1	A	329	LEU	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	873	18/19	0.96	0.21	-	35,37,43,43	0
3	BRU	T	840[B]	20/21	0.93	0.21	-	50,55,65,72	20
3	BRU	T	840[A]	20/21	0.93	0.21	-	42,47,68,127	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
4	DTP	A	875	30/30	0.82	0.25	0.60	62,73,91,92	0
5	MG	A	871	1/1	0.98	0.08	-	58,58,58,58	0

6.5 Other polymers (i)

There are no such residues in this entry.