



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 01:27 am BST

PDB ID : 2XGP
Title : Yeast DNA polymerase eta in complex with C8-2-acetylaminofluorene containing DNA
Authors : Scheider, S.; Lammens, K.; Schorr, S.; Hopfner, K.P.; Carell, T.
Deposited on : 2010-06-07
Resolution : 2.70 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 ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) 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	:	2.11
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)	:	2.11

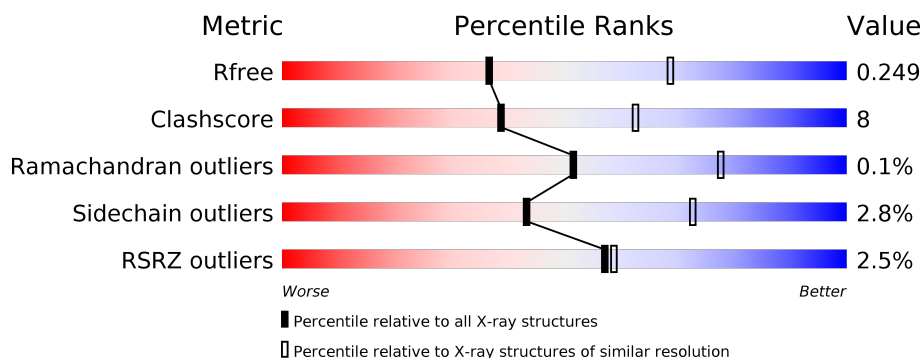
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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 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 $\leq 5\%$. 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	536	<div> <div>%</div> <div> <div></div> <div>78%</div> <div>16%</div> <div>• 5%</div> </div> </div>
1	B	536	<div> <div>4%</div> <div> <div></div> <div>77%</div> <div>18%</div> <div>• •</div> </div> </div>
2	P	9	<div> <div></div> <div> <div>56%</div> <div>44%</div> </div> </div>
2	Q	9	<div> <div>11%</div> <div> <div></div> <div>67%</div> <div>33%</div> </div> </div>
3	T	11	<div> <div></div> <div> <div>55%</div> <div>45%</div> </div> </div>
3	U	11	<div> <div></div> <div> <div>55%</div> <div>9%</div> <div>27%</div> <div>9%</div> </div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 9069 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 ETA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	511	Total	C	N	O	S	39	2	1
			4069	2593	687	765	24			
1	B	512	Total	C	N	O	S	73	3	1
			4079	2598	689	767	25			

There are 46 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-22	MET	-	expression tag	UNP Q04049
A	-21	ALA	-	expression tag	UNP Q04049
A	-20	SER	-	expression tag	UNP Q04049
A	-19	TRP	-	expression tag	UNP Q04049
A	-18	SER	-	expression tag	UNP Q04049
A	-17	HIS	-	expression tag	UNP Q04049
A	-16	PRO	-	expression tag	UNP Q04049
A	-15	GLN	-	expression tag	UNP Q04049
A	-14	PHE	-	expression tag	UNP Q04049
A	-13	GLU	-	expression tag	UNP Q04049
A	-12	LYS	-	expression tag	UNP Q04049
A	-11	GLY	-	expression tag	UNP Q04049
A	-10	ALA	-	expression tag	UNP Q04049
A	-9	SER	-	expression tag	UNP Q04049
A	-8	THR	-	expression tag	UNP Q04049
A	-7	SER	-	expression tag	UNP Q04049
A	-6	LEU	-	expression tag	UNP Q04049
A	-5	TYR	-	expression tag	UNP Q04049
A	-4	LYS	-	expression tag	UNP Q04049
A	-3	LYS	-	expression tag	UNP Q04049
A	-2	ALA	-	expression tag	UNP Q04049
A	-1	GLY	-	expression tag	UNP Q04049
A	0	ARG	-	expression tag	UNP Q04049
B	-22	MET	-	expression tag	UNP Q04049
B	-21	ALA	-	expression tag	UNP Q04049

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-20	SER	-	expression tag	UNP Q04049
B	-19	TRP	-	expression tag	UNP Q04049
B	-18	SER	-	expression tag	UNP Q04049
B	-17	HIS	-	expression tag	UNP Q04049
B	-16	PRO	-	expression tag	UNP Q04049
B	-15	GLN	-	expression tag	UNP Q04049
B	-14	PHE	-	expression tag	UNP Q04049
B	-13	GLU	-	expression tag	UNP Q04049
B	-12	LYS	-	expression tag	UNP Q04049
B	-11	GLY	-	expression tag	UNP Q04049
B	-10	ALA	-	expression tag	UNP Q04049
B	-9	SER	-	expression tag	UNP Q04049
B	-8	THR	-	expression tag	UNP Q04049
B	-7	SER	-	expression tag	UNP Q04049
B	-6	LEU	-	expression tag	UNP Q04049
B	-5	TYR	-	expression tag	UNP Q04049
B	-4	LYS	-	expression tag	UNP Q04049
B	-3	LYS	-	expression tag	UNP Q04049
B	-2	ALA	-	expression tag	UNP Q04049
B	-1	GLY	-	expression tag	UNP Q04049
B	0	ARG	-	expression tag	UNP Q04049

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	P	9	Total	C	N	O	P	0	0	0
			188	90	39	51	8			
2	Q	9	Total	C	N	O	P	0	0	0
			188	90	39	51	8			

- Molecule 3 is a DNA chain called 5'-D(*CP*8FG*CP*TP*CP*AP*TP*CP*CP*AP*C)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	T	11	Total	C	N	O	P	0	0	0
			232	119	38	65	10			
3	U	10	Total	C	N	O	P	0	0	0
			216	110	35	61	10			

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

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

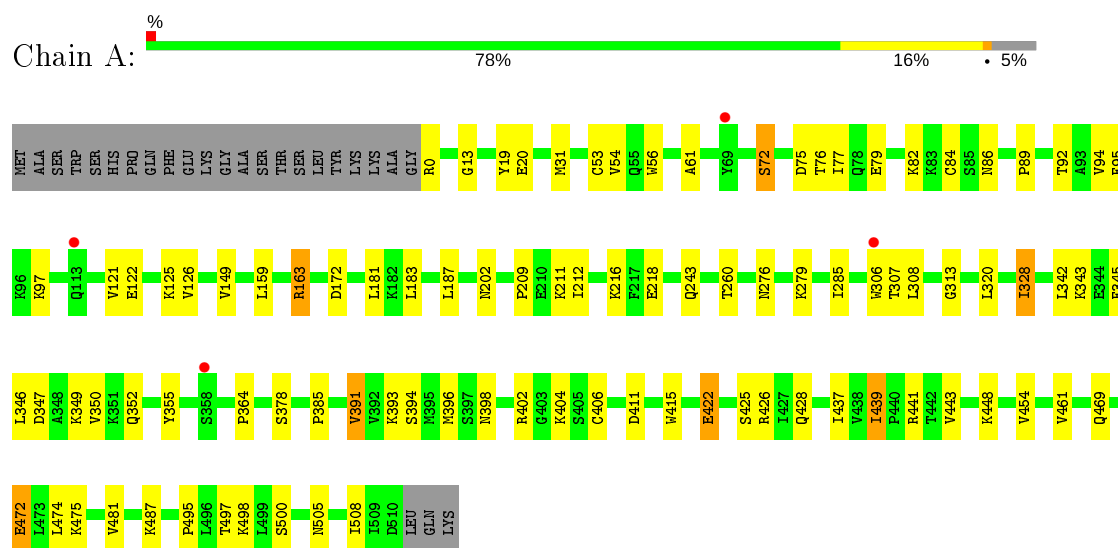
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	44	Total O 44 44	0	0
5	B	40	Total O 40 40	0	0
5	P	3	Total O 3 3	0	0
5	T	1	Total O 1 1	0	0
5	U	1	Total O 1 1	0	0

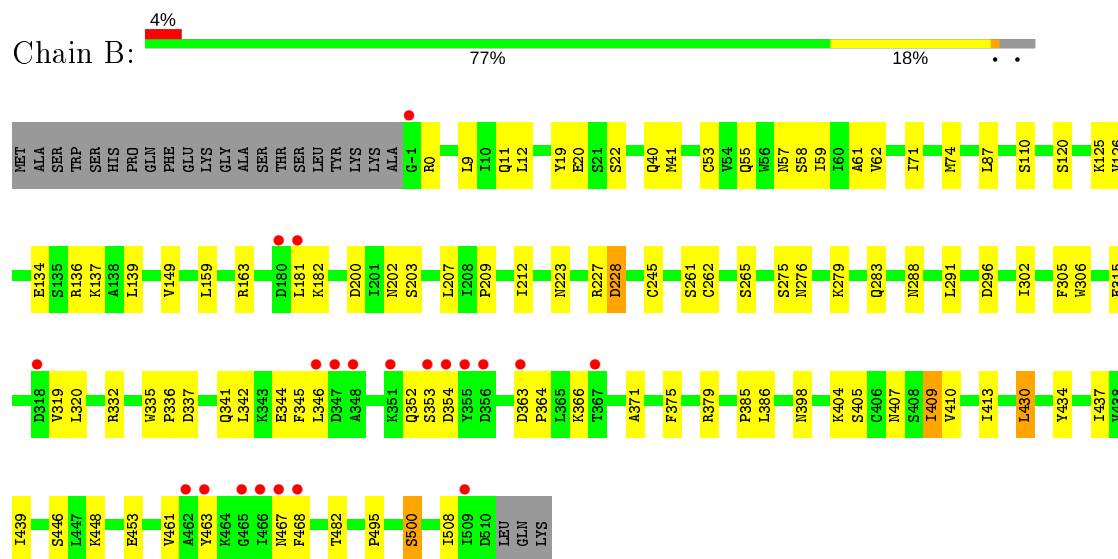
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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 ETA



• Molecule 1: DNA POLYMERASE ETA



• Molecule 2: 5'-D(*GP*TP*GP*GP*AP*TP*GP*AP*G)-3'





- Molecule 2: 5'-D(*GP*TP*GP*GP*AP*TP*GP*AP*G)-3'



- Molecule 3: 5'-D(*CP*8FG*CP*TP*CP*AP*TP*CP*CP*AP*C)-3'



- Molecule 3: 5'-D(*CP*8FG*CP*TP*CP*AP*TP*CP*CP*AP*C)-3'



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	103.44Å 103.44Å 292.35Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.76 – 2.70 48.76 – 2.69	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.76-2.70) 99.0 (48.76-2.69)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.76 (at 2.69Å)	Xtriage
Refinement program	REFMAC 5.5.0110	Depositor
R, R_{free}	0.220 , 0.272 0.210 , 0.249	Depositor DCC
R_{free} test set	3592 reflections (8.05%)	wwPDB-VP
Wilson B-factor (Å ²)	40.9	Xtriage
Anisotropy	0.033	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 34.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	9069	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.18% 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, 0.333 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: CA, 8FG

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	0.70	3/4153 (0.1%)	0.69	1/5600 (0.0%)
1	B	0.67	0/4166	0.67	0/5616
2	P	0.94	0/212	1.64	4/327 (1.2%)
2	Q	0.88	0/212	1.45	1/327 (0.3%)
3	T	0.96	0/213	1.63	4/322 (1.2%)
3	U	0.97	0/196	1.74	4/298 (1.3%)
All	All	0.71	3/9152 (0.0%)	0.83	14/12490 (0.1%)

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
1	B	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	475	LYS	CB-CG	-7.07	1.33	1.52
1	A	487	LYS	CD-CE	-6.42	1.35	1.51
1	A	82	LYS	CB-CG	5.51	1.67	1.52

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	U	6	DC	O4'-C4'-C3'	-10.67	99.60	106.00
2	P	7	DG	O4'-C1'-N9	-7.59	102.69	108.00
2	P	5	DG	O4'-C1'-N9	-7.53	102.73	108.00
3	U	8	DC	O4'-C1'-N1	5.88	112.12	108.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	T	4	DC	N1-C1'-C2'	5.79	123.61	112.60
3	T	12	DC	O4'-C1'-N1	5.72	112.01	108.00
3	T	14	DC	C1'-O4'-C4'	-5.68	104.42	110.10
1	A	0	ARG	CB-CG-CD	-5.62	96.98	111.60
3	U	7	DT	N3-C4-O4	5.57	123.24	119.90
2	P	11	DG	O4'-C1'-N9	5.44	111.81	108.00
3	U	8	DC	O4'-C1'-C2'	-5.16	101.77	105.90
2	P	8	DG	C5-C6-O6	-5.12	125.53	128.60
2	Q	7	DG	O4'-C1'-N9	-5.09	104.44	108.00
3	T	9	DA	O4'-C1'-N9	-5.04	104.47	108.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	74[B]	MET	Mainchain

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	A	4069	0	4126	60	0
1	B	4079	0	4135	65	0
2	P	188	0	101	0	0
2	Q	188	0	101	2	0
3	T	232	0	136	1	0
3	U	216	0	124	6	0
4	A	4	0	0	0	0
4	B	4	0	0	0	0
5	A	44	0	0	0	0
5	B	40	0	0	0	0
5	P	3	0	0	0	0
5	T	1	0	0	1	0
5	U	1	0	0	0	0
All	All	9069	0	8723	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 8.

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:345:PHE:O	1:A:349:LYS:HG2	1.82	0.80
3:U:5:8FG:H4'	3:U:6:DC:OP1	1.82	0.79
1:B:404:LYS:HG2	1:B:495:PRO:HG2	1.65	0.78
1:B:262:CYS:HB3	1:B:283:GLN:HG3	1.68	0.74
1:B:342:LEU:HD23	1:B:371:ALA:HA	1.68	0.73
1:B:404:LYS:HB3	1:B:407:ASN:OD1	1.88	0.73
1:A:306:TRP:O	1:A:307:THR:HG23	1.90	0.72
3:U:7:DT:H2'	3:U:8:DC:C6	2.26	0.71
1:B:398:ASN:HB3	1:B:500:SER:HB2	1.72	0.70
1:A:350:VAL:HG13	1:A:355:TYR:CE2	2.30	0.67
1:A:31:MET:HG2	1:A:260:THR:HG22	1.77	0.67
1:B:136:ARG:NH1	1:B:139:LEU:HD23	2.10	0.65
1:B:398:ASN:CB	1:B:500:SER:HB2	2.25	0.65
1:A:149:VAL:HG22	1:A:159:LEU:HD22	1.80	0.63
1:A:13:GLY:O	1:A:163[A]:ARG:HD2	1.99	0.62
1:A:398:ASN:HD22	1:A:500:SER:HB3	1.64	0.61
1:B:363:ASP:HB3	1:B:366:LYS:HB2	1.81	0.61
1:A:306:TRP:C	1:A:307:THR:HG23	2.21	0.61
1:B:228:ASP:OD2	1:B:228:ASP:N	2.34	0.61
3:U:5:8FG:C4'	3:U:6:DC:OP1	2.49	0.60
1:A:328[A]:ILE:CD1	1:A:378:SER:HA	2.32	0.60
1:A:125:LYS:NZ	1:A:398:ASN:O	2.29	0.58
1:B:336:PRO:HD2	1:B:341:GLN:OE1	2.04	0.57
1:A:404:LYS:HD3	1:A:495:PRO:HG3	1.87	0.56
1:A:19:TYR:CE2	1:A:385:PRO:HB3	2.41	0.56
1:A:393:LYS:NZ	3:T:11:DC:OP2	2.26	0.55
1:B:209:PRO:O	1:B:212:ILE:HG22	2.08	0.54
3:U:7:DT:H2''	3:U:8:DC:O4'	2.08	0.53
1:A:396:MET:O	1:A:426:ARG:NH2	2.41	0.53
1:B:315:GLU:O	1:B:319:VAL:HG23	2.07	0.53
1:A:183:LEU:HD22	1:A:187:LEU:HD12	1.90	0.53
3:U:5:8FG:H2''	3:U:6:DC:O5'	2.09	0.52
1:A:306:TRP:O	1:A:307:THR:CG2	2.57	0.52
1:A:97:LYS:HG3	1:A:415:TRP:CE2	2.44	0.52
1:B:149:VAL:HG22	1:B:159:LEU:HD22	1.92	0.52
1:B:352:GLN:HB3	1:B:354:ASP:OD1	2.10	0.52
1:B:467:ASN:OD1	1:B:467:ASN:C	2.48	0.52
1:B:363:ASP:OD2	1:B:364:PRO:HD2	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:125:LYS:HE2	1:B:398:ASN:O	2.10	0.51
1:A:306:TRP:O	1:A:307:THR:OG1	2.21	0.51
1:A:218:GLU:O	1:A:285:ILE:HD11	2.11	0.51
1:B:40:GLN:NE2	1:B:41:MET:CE	2.74	0.51
1:B:59:ILE:HG21	1:B:62:VAL:HG22	1.93	0.51
1:A:76:THR:OG1	1:A:79:GLU:HG3	2.11	0.51
1:A:183:LEU:HD22	1:A:187:LEU:CD1	2.40	0.50
1:A:306:TRP:O	1:A:307:THR:CB	2.59	0.50
1:B:409:ILE:O	1:B:413:ILE:HG13	2.10	0.50
1:B:439:ILE:HD11	1:B:463:TYR:CD2	2.46	0.50
1:B:337:ASP:HB2	1:B:341:GLN:NE2	2.27	0.50
1:A:343:LYS:NZ	1:A:347:ASP:OD2	2.45	0.50
1:B:245:CYS:SG	1:B:262:CYS:SG	3.09	0.50
1:B:306:TRP:CZ3	2:Q:12:DA:H4'	2.47	0.50
1:A:443:VAL:HG12	1:A:461:VAL:CG2	2.42	0.49
1:B:448:LYS:HA	1:B:453:GLU:O	2.13	0.49
1:A:122:GLU:OE2	1:A:402:ARG:O	2.30	0.48
1:B:409:ILE:HG22	1:B:410:VAL:N	2.28	0.48
1:B:134:GLU:OE2	1:B:137:LYS:HE3	2.13	0.48
1:A:391:VAL:HG21	1:A:508:ILE:HD12	1.95	0.47
1:B:296:ASP:OD2	1:B:332:ARG:NH2	2.46	0.47
1:B:200:ASP:O	1:B:203:SER:HB2	2.14	0.47
1:A:437:ILE:HG22	1:A:439:ILE:HD11	1.97	0.47
1:B:320:LEU:HD21	1:B:346:LEU:HD21	1.97	0.47
1:A:441:ARG:O	1:A:461:VAL:N	2.40	0.47
1:B:409:ILE:CG2	1:B:410:VAL:N	2.78	0.46
1:B:404:LYS:O	1:B:407:ASN:OD1	2.34	0.46
1:A:54:VAL:O	1:A:92:THR:HG23	2.15	0.46
2:Q:11:DG:C2'	2:Q:12:DA:H5'	2.46	0.46
1:A:172:ASP:C	1:A:172:ASP:OD1	2.54	0.45
1:B:55:GLN:HG3	1:B:126:VAL:HG21	1.98	0.45
1:B:136:ARG:HH12	1:B:139:LEU:HD23	1.81	0.45
1:B:439:ILE:HG13	1:B:468:PHE:HB2	1.99	0.45
1:B:53:CYS:HB3	1:B:61:ALA:HB3	1.98	0.45
1:A:350:VAL:HG11	1:A:364:PRO:HB3	1.98	0.45
1:A:94:VAL:HG12	1:A:126:VAL:HA	1.97	0.45
1:A:216:LYS:O	1:A:243:GLN:NE2	2.48	0.45
1:A:391:VAL:CG2	1:A:508:ILE:HD12	2.46	0.45
1:B:57:ASN:ND2	1:B:110:SER:OG	2.50	0.45
1:B:181:LEU:HD12	1:B:182:LYS:H	1.82	0.44
1:B:375:PHE:CE1	1:B:379:ARG:NH2	2.85	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:437:ILE:CG2	1:A:439:ILE:HD11	2.46	0.44
1:B:320:LEU:HD21	1:B:346:LEU:CD2	2.47	0.44
1:A:84:CYS:O	1:A:86:ASN:N	2.48	0.44
1:A:95:PHE:CE1	1:A:422:GLU:HG3	2.53	0.44
1:B:409:ILE:HD11	1:B:482:THR:HA	1.99	0.44
1:B:386:LEU:HD12	1:B:386:LEU:N	2.32	0.44
3:U:6:DC:H2'	3:U:7:DT:C6	2.52	0.44
1:A:72:SER:O	1:A:75:ASP:HB2	2.18	0.43
1:B:430:LEU:HD23	1:B:430:LEU:HA	1.76	0.43
1:A:448:LYS:HG3	1:A:454:VAL:HG23	2.00	0.43
1:B:0:ARG:O	1:B:207:LEU:HD13	2.18	0.43
1:B:342:LEU:HD23	1:B:371:ALA:CA	2.44	0.43
1:A:308:LEU:O	1:A:313:GLY:HA3	2.19	0.43
1:B:212:ILE:HA	1:B:212:ILE:HD12	1.86	0.43
1:B:288:ASN:HD22	1:B:291:LEU:HD11	1.83	0.43
1:B:276:ASN:OD1	1:B:279:LYS:NZ	2.36	0.43
1:B:344:GLU:O	1:B:345:PHE:C	2.57	0.43
1:B:354:ASP:OD1	1:B:354:ASP:N	2.48	0.43
1:A:209:PRO:O	1:A:212:ILE:HG22	2.19	0.43
1:A:181:LEU:HD21	1:A:211:LYS:HB3	2.01	0.43
1:A:77:ILE:HD11	1:A:89:PRO:HB3	2.01	0.43
1:A:437:ILE:HD11	1:B:202:ASN:ND2	2.34	0.42
1:A:20:GLU:O	1:A:20:GLU:HG2	2.18	0.42
1:A:342:LEU:HD12	1:A:342:LEU:O	2.19	0.42
1:A:394:SER:HB3	5:T:2001:HOH:O	2.20	0.42
1:A:469:GLN:HB3	1:A:472:GLU:HB2	2.02	0.42
1:B:19:TYR:CE2	1:B:385:PRO:HB3	2.55	0.42
1:A:56:TRP:CE3	1:A:121:VAL:HG22	2.54	0.42
1:A:276:ASN:O	1:A:279:LYS:HD2	2.20	0.42
1:A:306:TRP:CE3	1:A:307:THR:CG2	3.03	0.42
1:A:320:LEU:HD21	1:A:346:LEU:HD21	2.01	0.42
1:B:11:GLN:NE2	1:B:20:GLU:HB2	2.34	0.42
1:A:202:ASN:OD1	1:B:437:ILE:HD11	2.20	0.42
1:B:261:SER:HB3	1:B:275:SER:O	2.20	0.41
1:A:441:ARG:HB2	1:A:505:ASN:HB3	2.01	0.41
1:B:335:TRP:CE2	1:B:342:LEU:HD12	2.55	0.41
1:B:181:LEU:HD12	1:B:181:LEU:HA	1.83	0.41
1:B:9:LEU:O	1:B:12:LEU:HB2	2.20	0.41
1:B:71:ILE:CD1	1:B:87:LEU:HD13	2.51	0.41
1:A:448:LYS:HB2	1:A:498:LYS:HB2	2.02	0.41
1:B:302:ILE:HD12	1:B:305:PHE:CE1	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:59:ILE:CG2	1:B:62:VAL:HG22	2.50	0.41
1:A:307:THR:O	1:A:308:LEU:HD23	2.21	0.41
1:A:328[A]:ILE:H	1:A:328[A]:ILE:HG22	1.54	0.41
1:A:391:VAL:CG2	1:A:508:ILE:CD1	2.98	0.41
1:A:53:CYS:HB3	1:A:61:ALA:HB3	2.03	0.41
1:A:406:CYS:HA	1:A:411:ASP:HB3	2.02	0.40
1:B:345:PHE:CD2	1:B:345:PHE:C	2.95	0.40
1:B:363:ASP:HA	1:B:364:PRO:HD3	1.86	0.40
1:B:223:ASN:ND2	1:B:227:ARG:O	2.50	0.40
1:B:430:LEU:HD13	1:B:434:TYR:CE2	2.57	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	511/536 (95%)	497 (97%)	14 (3%)	0	100	100
1	B	513/536 (96%)	499 (97%)	13 (2%)	1 (0%)	47	73
All	All	1024/1072 (96%)	996 (97%)	27 (3%)	1 (0%)	51	78

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	508	ILE

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	459/478 (96%)	444 (97%)	15 (3%)	38	67
1	B	460/478 (96%)	447 (97%)	13 (3%)	43	73
All	All	919/956 (96%)	891 (97%)	28 (3%)	43	70

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	72	SER
1	A	163[A]	ARG
1	A	163[B]	ARG
1	A	328[A]	ILE
1	A	328[B]	ILE
1	A	352	GLN
1	A	391	VAL
1	A	422	GLU
1	A	425	SER
1	A	428	GLN
1	A	439	ILE
1	A	472	GLU
1	A	474	LEU
1	A	481	VAL
1	A	497	THR
1	B	22	SER
1	B	58	SER
1	B	120	SER
1	B	163	ARG
1	B	228	ASP
1	B	265	SER
1	B	353	SER
1	B	405	SER
1	B	409	ILE
1	B	430	LEU
1	B	446	SER
1	B	461	VAL
1	B	500	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	57	ASN
1	A	288	ASN
1	A	398	ASN
1	A	400	ASN
1	B	32	ASN
1	B	40	GLN
1	B	57	ASN
1	B	288	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 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
3	8FG	T	5	3	38,44,45	1.48	6 (15%)	49,66,69	1.65	7 (14%)
3	8FG	U	5	3	38,44,45	1.43	6 (15%)	49,66,69	1.55	6 (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
3	8FG	T	5	3	-	2/11/41/42	0/6/6/6
3	8FG	U	5	3	-	4/11/41/42	0/6/6/6

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	U	5	8FG	C6-C5	4.56	1.49	1.41
3	U	5	8FG	C8-N29	4.39	1.46	1.35
3	T	5	8FG	C8-N29	4.19	1.45	1.35
3	T	5	8FG	C6-C5	3.79	1.47	1.41
3	T	5	8FG	C39-C35	-3.63	1.37	1.46
3	U	5	8FG	C37-C38	-2.79	1.44	1.51
3	U	5	8FG	C39-C35	-2.78	1.39	1.46
3	U	5	8FG	C37-C34	-2.77	1.44	1.51
3	T	5	8FG	C37-C38	-2.76	1.44	1.51
3	T	5	8FG	C32-N29	-2.73	1.40	1.44
3	T	5	8FG	C37-C34	-2.64	1.44	1.51
3	U	5	8FG	C32-N29	-2.25	1.41	1.44

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	U	5	8FG	C2-N3-C4	5.99	122.19	115.36
3	T	5	8FG	C2'-C1'-N9	-5.49	110.01	116.01
3	T	5	8FG	C2-N3-C4	5.05	121.12	115.36
3	U	5	8FG	N3-C2-N1	-4.13	121.72	127.22
3	T	5	8FG	C5-C6-N1	-3.78	118.27	123.43
3	T	5	8FG	N3-C2-N1	-3.63	122.38	127.22
3	U	5	8FG	C5-C6-N1	-3.49	118.66	123.43
3	U	5	8FG	C6-N1-C2	2.91	120.55	115.93
3	T	5	8FG	C6-N1-C2	2.90	120.53	115.93
3	T	5	8FG	C6-C5-C4	-2.81	118.11	120.80
3	U	5	8FG	C6-C5-C4	-2.65	118.27	120.80
3	T	5	8FG	C32-N29-C30	-2.22	117.13	121.52
3	U	5	8FG	C38-C37-C34	2.03	106.88	102.73

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	U	5	8FG	O44-C30-N29-C8
3	U	5	8FG	C45-C30-N29-C8
3	U	5	8FG	O44-C30-N29-C32
3	U	5	8FG	C45-C30-N29-C32
3	T	5	8FG	C3'-C4'-C5'-O5'
3	T	5	8FG	O4'-C4'-C5'-O5'

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	U	5	8FG	3	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 8 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	511/536 (95%)	-0.07	4 (0%) 86 87	9, 26, 48, 58	10 (1%)
1	B	512/536 (95%)	0.02	21 (4%) 37 36	13, 26, 61, 74	22 (4%)
2	P	9/9 (100%)	-0.33	0 100 100	27, 32, 40, 40	0
2	Q	9/9 (100%)	0.13	1 (11%) 5 4	32, 39, 87, 95	0
3	T	10/11 (90%)	-0.26	0 100 100	21, 36, 44, 68	0
3	U	9/11 (81%)	0.15	0 100 100	36, 39, 78, 88	0
All	All	1060/1112 (95%)	-0.02	26 (2%) 57 59	9, 26, 56, 95	32 (3%)

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	355	TYR	4.4
1	B	318	ASP	3.6
1	B	346	LEU	3.6
1	B	465	GLY	3.5
1	B	180	ASP	3.4
1	B	356	ASP	3.4
1	A	113	GLN	3.0
1	B	-1	GLY	3.0
1	B	353	SER	3.0
1	B	466	ILE	2.9
1	B	363	ASP	2.9
1	B	463	TYR	2.8
1	B	468	PHE	2.7
1	B	467	ASN	2.6
1	B	348	ALA	2.6
1	B	347	ASP	2.5
1	B	181	LEU	2.5
1	A	69	TYR	2.4
1	B	354	ASP	2.4

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Mol	Chain	Res	Type	RSRZ
2	Q	13	DG	2.3
1	B	367	THR	2.3
1	A	306	TRP	2.2
1	B	462	ALA	2.2
1	B	351	LYS	2.2
1	A	358	SER	2.1
1	B	509	ILE	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. 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	B-factors(Å ²)	Q<0.9
3	8FG	U	5	39/40	0.74	0.31	86,91,103,104	0
3	8FG	T	5	39/40	0.94	0.15	23,35,59,60	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. 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	B-factors(Å ²)	Q<0.9
4	CA	B	1511	1/1	0.69	0.24	69,69,69,69	0
4	CA	B	1512	1/1	0.75	0.17	60,60,60,60	0
4	CA	A	1513	1/1	0.90	0.08	50,50,50,50	0
4	CA	B	1513	1/1	0.91	0.21	59,59,59,59	0
4	CA	A	1512	1/1	0.91	0.08	70,70,70,70	0
4	CA	B	1514	1/1	0.92	0.07	53,53,53,53	0
4	CA	A	1511	1/1	0.92	0.14	60,60,60,60	0
4	CA	A	1514	1/1	0.95	0.18	3,3,3,3	1

6.5 Other polymers [i](#)

There are no such residues in this entry.