



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 11:47 AM GMT

PDB ID : 3PNS
Title : Crystal Structure of Uridine Phosphorylase Complexed with Uracil from *Vibrio cholerae* O1 biovar El Tor
Authors : Maltseva, N.; Kim, Y.; Hasseman, J.; Anderson, W.F.; Joachimiak, A.; Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on : 2010-11-19
Resolution : 2.00 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

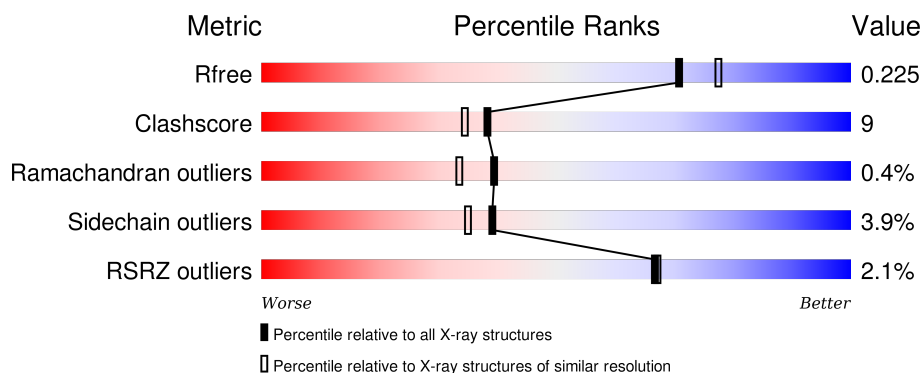
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	91344	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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 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	261	<div> <div>2%</div> <div> <div></div> <div>77%</div> <div>17%</div> <div>• •</div> </div> </div>
1	B	261	<div> <div>%</div> <div> <div></div> <div>77%</div> <div>15%</div> <div>• 7%</div> </div> </div>
1	C	261	<div> <div>2%</div> <div> <div></div> <div>70%</div> <div>22%</div> <div>• 5%</div> </div> </div>
1	D	261	<div> <div>%</div> <div> <div></div> <div>78%</div> <div>14%</div> <div>• 6%</div> </div> </div>
1	E	261	<div> <div>3%</div> <div> <div></div> <div>75%</div> <div>18%</div> <div>• 6%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	F	261	
1	G	261	
1	H	261	
1	I	261	
1	J	261	
1	K	261	
1	L	261	

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	GOL	C	275	-	-	X	-
2	GOL	D	271	-	-	-	X
2	GOL	F	270	-	-	-	X
2	GOL	G	271	-	-	-	X
2	GOL	I	272	-	-	-	X
2	GOL	K	271	-	-	-	X
3	URA	B	271	-	-	-	X
3	URA	C	276	-	-	-	X
6	FMT	E	273	-	-	X	-
6	FMT	H	272	-	-	X	-

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 25215 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 Uridine phosphorylase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	250	Total	C	N	O	S	Se	0	9	0
			1951	1216	346	375	4	10			
1	B	244	Total	C	N	O	S	Se	0	10	0
			1902	1183	334	371	3	11			
1	C	248	Total	C	N	O	S	Se	0	8	0
			1925	1199	341	372	3	10			
1	D	245	Total	C	N	O	S	Se	0	8	0
			1896	1183	337	363	3	10			
1	E	245	Total	C	N	O	S	Se	0	4	0
			1864	1165	327	359	3	10			
1	F	250	Total	C	N	O	S	Se	1	7	0
			1923	1202	337	370	4	10			
1	G	250	Total	C	N	O	S	Se	0	6	0
			1916	1197	334	371	3	11			
1	H	249	Total	C	N	O	S	Se	1	15	0
			1985	1235	353	383	4	10			
1	I	250	Total	C	N	O	S	Se	0	9	0
			1937	1211	334	378	4	10			
1	J	241	Total	C	N	O	S	Se	1	4	0
			1826	1142	316	354	4	10			
1	K	251	Total	C	N	O	S	Se	0	9	0
			1941	1211	335	381	4	10			
1	L	245	Total	C	N	O	S	Se	0	6	0
			1879	1173	329	364	3	10			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	EXPRESSION TAG	UNP Q9KT71
A	-1	ASN	-	EXPRESSION TAG	UNP Q9KT71
A	0	ALA	-	EXPRESSION TAG	UNP Q9KT71
B	-2	SER	-	EXPRESSION TAG	UNP Q9KT71
B	-1	ASN	-	EXPRESSION TAG	UNP Q9KT71

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Chain	Residue	Modelled	Actual	Comment	Reference
B	0	ALA	-	EXPRESSION TAG	UNP Q9KT71
C	-2	SER	-	EXPRESSION TAG	UNP Q9KT71
C	-1	ASN	-	EXPRESSION TAG	UNP Q9KT71
C	0	ALA	-	EXPRESSION TAG	UNP Q9KT71
D	-2	SER	-	EXPRESSION TAG	UNP Q9KT71
D	-1	ASN	-	EXPRESSION TAG	UNP Q9KT71
D	0	ALA	-	EXPRESSION TAG	UNP Q9KT71
E	-2	SER	-	EXPRESSION TAG	UNP Q9KT71
E	-1	ASN	-	EXPRESSION TAG	UNP Q9KT71
E	0	ALA	-	EXPRESSION TAG	UNP Q9KT71
F	-2	SER	-	EXPRESSION TAG	UNP Q9KT71
F	-1	ASN	-	EXPRESSION TAG	UNP Q9KT71
F	0	ALA	-	EXPRESSION TAG	UNP Q9KT71
G	-2	SER	-	EXPRESSION TAG	UNP Q9KT71
G	-1	ASN	-	EXPRESSION TAG	UNP Q9KT71
G	0	ALA	-	EXPRESSION TAG	UNP Q9KT71
H	-2	SER	-	EXPRESSION TAG	UNP Q9KT71
H	-1	ASN	-	EXPRESSION TAG	UNP Q9KT71
H	0	ALA	-	EXPRESSION TAG	UNP Q9KT71
I	-2	SER	-	EXPRESSION TAG	UNP Q9KT71
I	-1	ASN	-	EXPRESSION TAG	UNP Q9KT71
I	0	ALA	-	EXPRESSION TAG	UNP Q9KT71
J	-2	SER	-	EXPRESSION TAG	UNP Q9KT71
J	-1	ASN	-	EXPRESSION TAG	UNP Q9KT71
J	0	ALA	-	EXPRESSION TAG	UNP Q9KT71
K	-2	SER	-	EXPRESSION TAG	UNP Q9KT71
K	-1	ASN	-	EXPRESSION TAG	UNP Q9KT71
K	0	ALA	-	EXPRESSION TAG	UNP Q9KT71
L	-2	SER	-	EXPRESSION TAG	UNP Q9KT71
L	-1	ASN	-	EXPRESSION TAG	UNP Q9KT71
L	0	ALA	-	EXPRESSION TAG	UNP Q9KT71

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



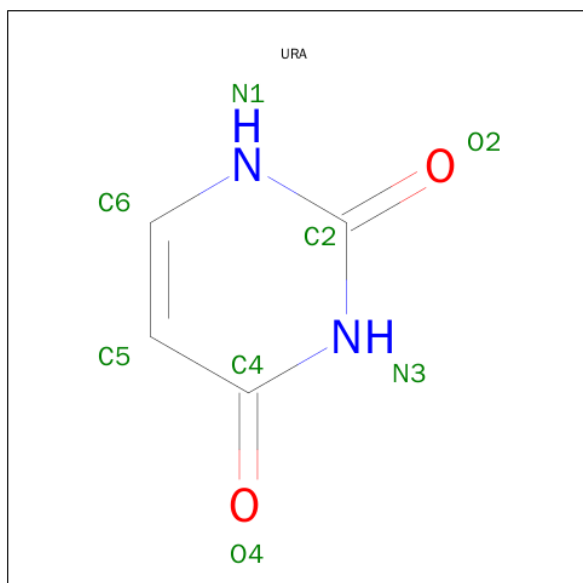
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		
2	C	1	Total	C	O	0	0
			6	3	3		
2	C	1	Total	C	O	0	0
			6	3	3		
2	C	1	Total	C	O	0	0
			6	3	3		
2	D	1	Total	C	O	0	0
			6	3	3		
2	D	1	Total	C	O	0	0
			6	3	3		
2	E	1	Total	C	O	0	0
			6	3	3		
2	E	1	Total	C	O	0	0
			6	3	3		
2	F	1	Total	C	O	0	0
			6	3	3		
2	F	1	Total	C	O	0	0
			6	3	3		
2	G	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	G	1	Total	C	O	0	0
			6	3	3		
2	G	1	Total	C	O	0	0
			6	3	3		
2	H	1	Total	C	O	0	0
			6	3	3		
2	I	1	Total	C	O	0	0
			6	3	3		
2	I	1	Total	C	O	0	0
			6	3	3		
2	J	1	Total	C	O	0	0
			6	3	3		
2	J	1	Total	C	O	0	0
			6	3	3		
2	K	1	Total	C	O	0	0
			6	3	3		
2	K	1	Total	C	O	0	0
			6	3	3		
2	L	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is URACIL (three-letter code: URA) (formula: $C_4H_4N_2O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	N	O	0	0
			8	4	2	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	C	1	Total 8	C 4	N 2	O 2	0	0
3	D	1	Total 8	C 4	N 2	O 2	0	0
3	E	1	Total 8	C 4	N 2	O 2	0	0
3	F	1	Total 8	C 4	N 2	O 2	0	0
3	G	1	Total 8	C 4	N 2	O 2	0	0
3	H	1	Total 8	C 4	N 2	O 2	0	0
3	I	1	Total 8	C 4	N 2	O 2	0	0
3	J	1	Total 8	C 4	N 2	O 2	0	0
3	K	1	Total 8	C 4	N 2	O 2	0	0
3	L	1	Total 8	C 4	N 2	O 2	0	0

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

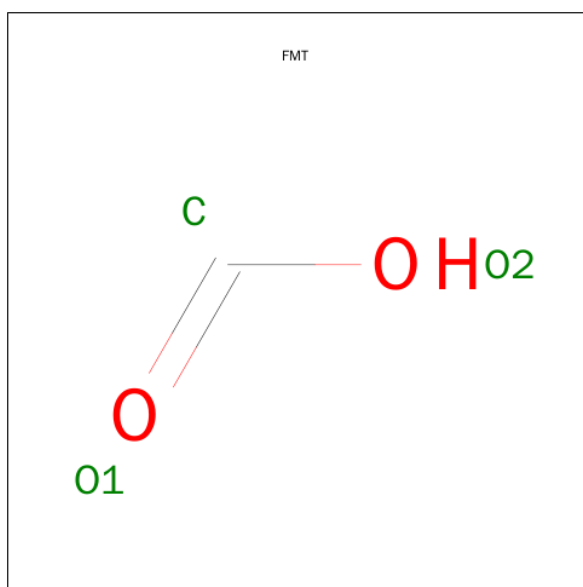
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	I	2	Total 2	Cl 2	0	0
4	C	1	Total 1	Cl 1	0	0
4	E	1	Total 1	Cl 1	0	0

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	C	1	Total	O	S	0	0
			5	4	1		
5	G	1	Total	O	S	0	0
			5	4	1		
5	I	1	Total	O	S	0	0
			5	4	1		

- Molecule 6 is FORMIC ACID (three-letter code: FMT) (formula: CH_2O_2).



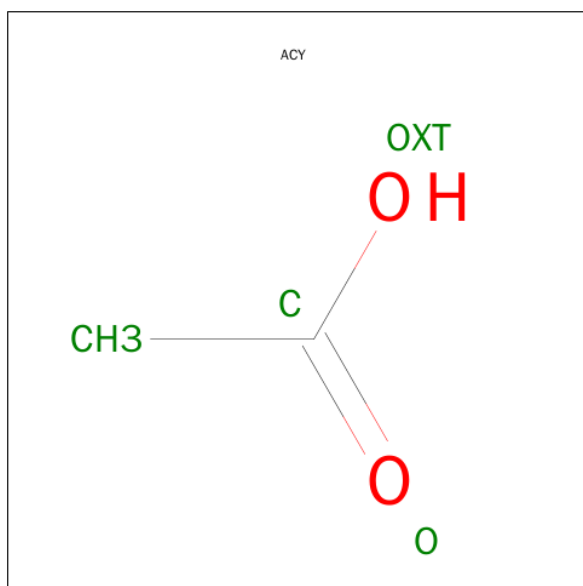
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	E	1	Total	C	O	0	0
			3	1	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	H	1	Total	C	O	0	0
			3	1	2		

- Molecule 7 is ACETIC ACID (three-letter code: ACY) (formula: C₂H₄O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	F	1	Total	C	O	0	0
			4	2	2		

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	161	Total	O	0	0
			161	161		
8	B	190	Total	O	0	0
			190	190		
8	C	132	Total	O	0	0
			132	132		
8	D	197	Total	O	0	0
			197	197		
8	E	151	Total	O	0	0
			151	151		
8	F	161	Total	O	0	0
			161	161		
8	G	165	Total	O	0	0
			165	165		

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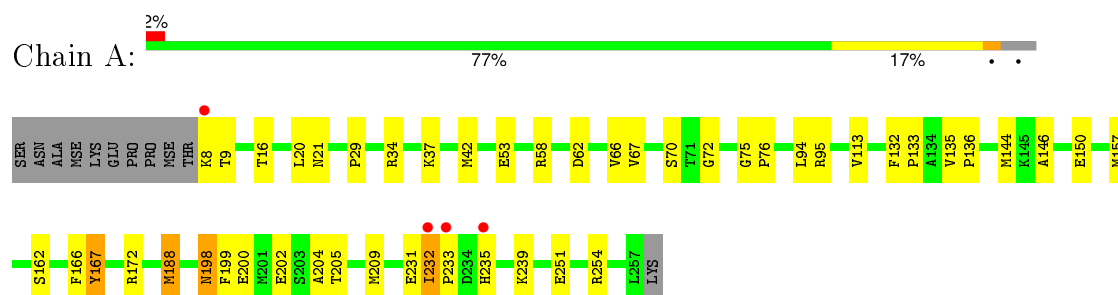
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	H	173	Total 173	O 173	0	0
8	I	181	Total 181	O 181	0	0
8	J	156	Total 156	O 156	0	0
8	K	171	Total 171	O 171	0	0
8	L	171	Total 171	O 171	0	0

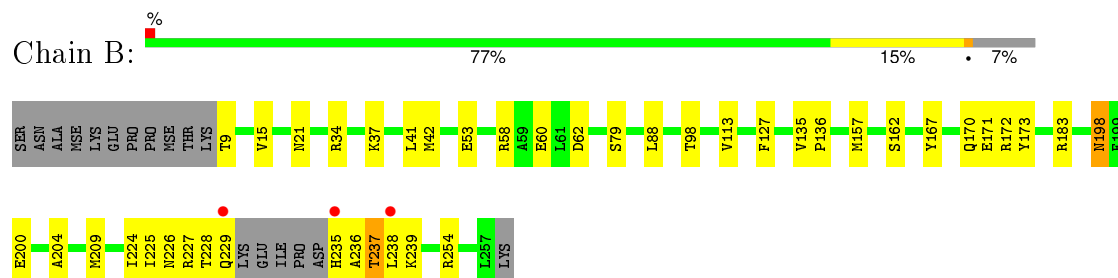
3 Residue-property plots

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

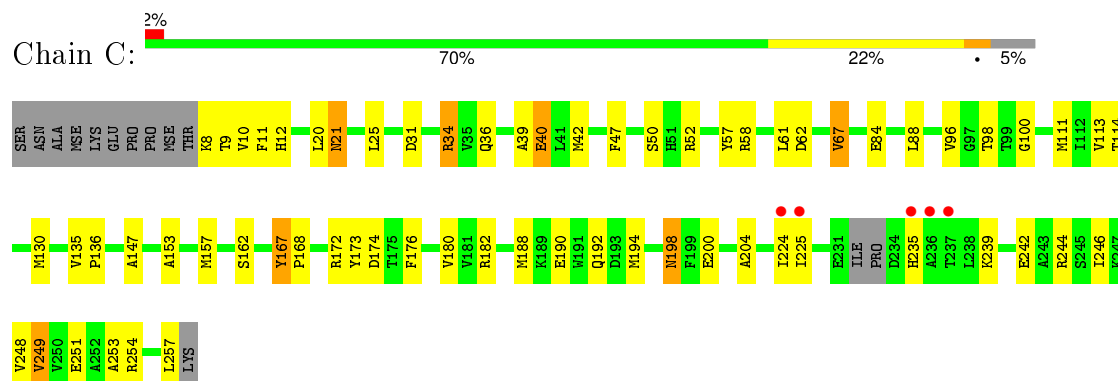
- Molecule 1: Uridine phosphorylase



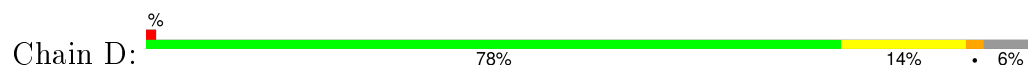
- Molecule 1: Uridine phosphorylase

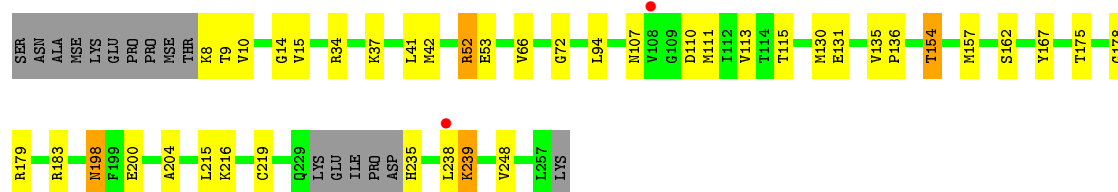


- Molecule 1: Uridine phosphorylase

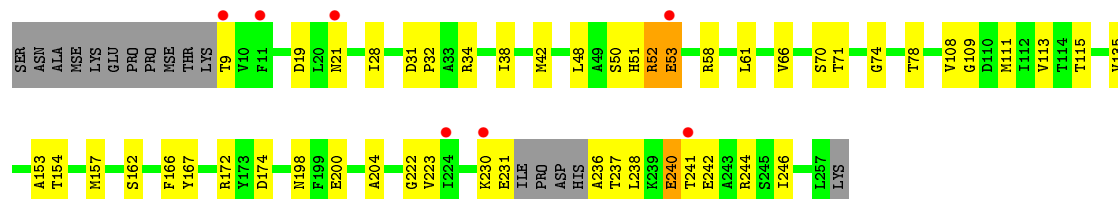
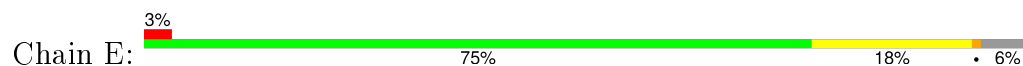


- Molecule 1: Uridine phosphorylase

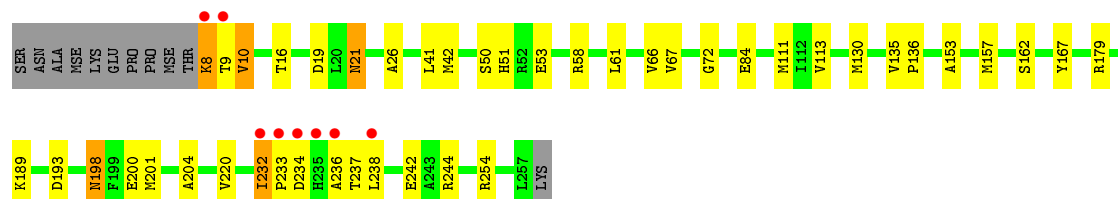
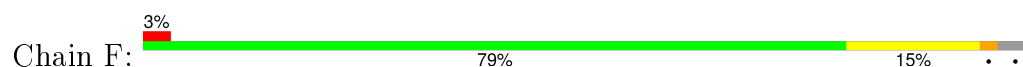




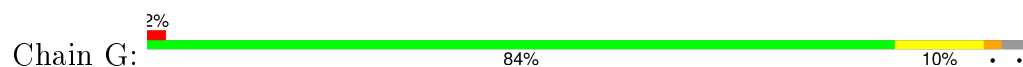
- Molecule 1: Uridine phosphorylase



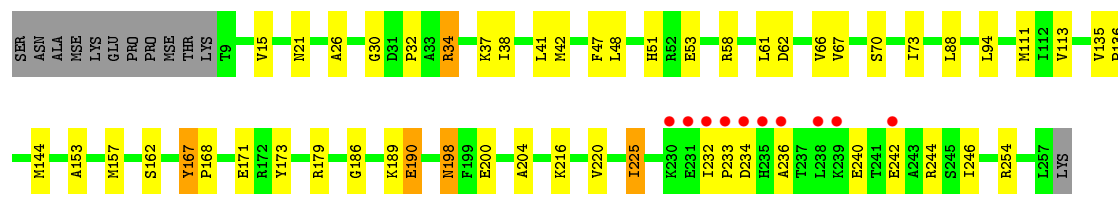
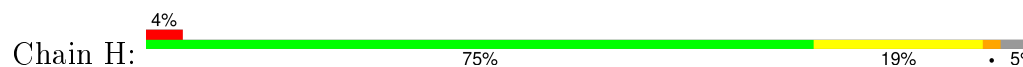
- Molecule 1: Uridine phosphorylase



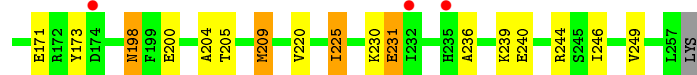
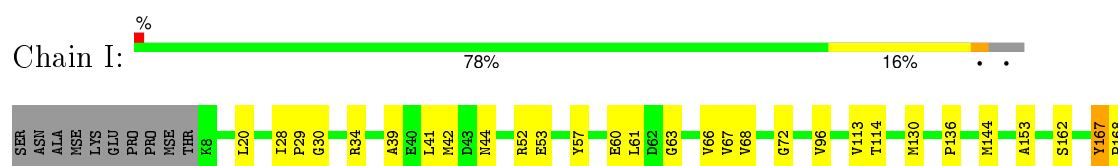
- Molecule 1: Uridine phosphorylase



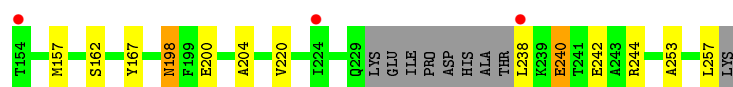
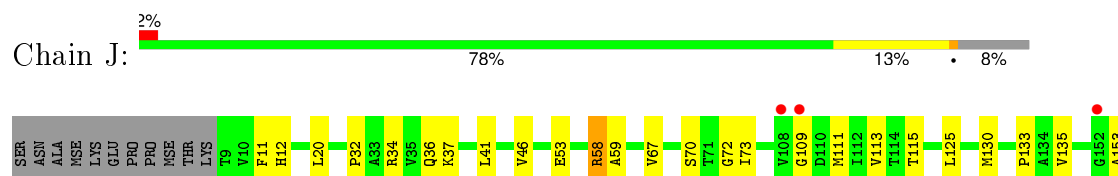
- Molecule 1: Uridine phosphorylase



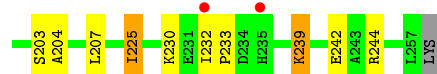
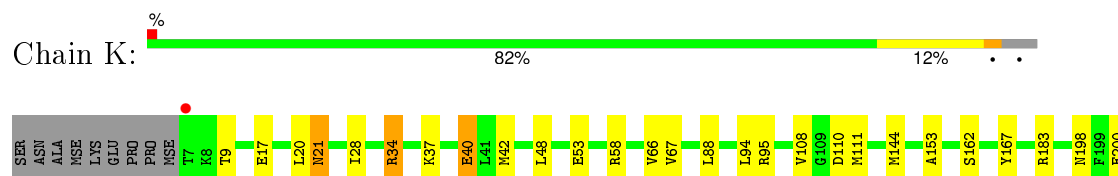
- Molecule 1: Uridine phosphorylase



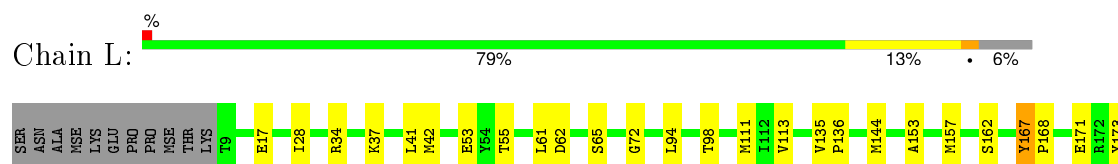
- Molecule 1: Uridine phosphorylase



- Molecule 1: Uridine phosphorylase



- Molecule 1: Uridine phosphorylase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	103.28Å 174.46Å 180.02Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.64 – 2.00 49.64 – 2.00	Depositor EDS
% Data completeness (in resolution range)	92.0 (49.64-2.00) 92.1 (49.64-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.09 (at 2.00Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.4_486)	Depositor
R, R_{free}	0.174 , 0.226 0.174 , 0.225	Depositor DCC
R_{free} test set	10079 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	22.8	Xtriage
Anisotropy	0.017	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 62.1	EDS
Estimated twinning fraction	0.000 for -h,l,k	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	6 of 200935 reflections (0.003%)	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	25215	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 43.31 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 1.7970e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: GOL, URA, CL, FMT, SO4, ACY

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	0.52	0/1972	0.65	0/2654
1	B	0.54	0/1921	0.66	0/2585
1	C	0.49	0/1944	0.61	0/2614
1	D	0.51	0/1915	0.62	0/2578
1	E	0.48	0/1882	0.61	0/2534
1	F	0.50	0/1945	0.63	0/2621
1	G	0.52	0/1937	0.63	1/2610 (0.0%)
1	H	0.50	0/2007	0.62	0/2703
1	I	0.50	0/1958	0.62	0/2640
1	J	0.51	0/1844	0.60	0/2484
1	K	0.53	0/1962	0.63	0/2645
1	L	0.55	0/1897	0.63	0/2553
All	All	0.51	0/23184	0.63	1/31221 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	182	ARG	NE-CZ-NH1	5.37	122.99	120.30

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	1951	0	1944	42	0
1	B	1902	0	1877	39	0
1	C	1925	0	1915	52	0
1	D	1896	0	1894	37	0
1	E	1864	0	1866	43	0
1	F	1923	0	1916	32	0
1	G	1916	0	1910	26	0
1	H	1985	0	1970	40	0
1	I	1937	0	1925	47	0
1	J	1826	0	1818	31	0
1	K	1941	0	1924	22	0
1	L	1879	0	1874	22	0
2	A	12	0	16	0	0
2	B	12	0	16	1	0
2	C	18	0	24	9	0
2	D	12	0	16	2	0
2	E	12	0	16	1	0
2	F	12	0	16	5	0
2	G	18	0	24	0	0
2	H	6	0	8	1	0
2	I	12	0	16	0	0
2	J	12	0	16	1	0
2	K	12	0	16	3	0
2	L	6	0	8	0	0
3	B	8	0	3	2	0
3	C	8	0	3	1	0
3	D	8	0	3	0	0
3	E	8	0	3	1	0
3	F	8	0	3	2	0
3	G	8	0	3	0	0
3	H	8	0	3	0	0
3	I	8	0	3	3	0
3	J	8	0	3	0	0
3	K	8	0	3	2	0
3	L	8	0	3	0	0
4	C	1	0	0	0	0
4	E	1	0	0	0	0
4	I	2	0	0	0	0
5	C	5	0	0	1	0
5	G	5	0	0	1	0
5	I	5	0	0	0	0
6	E	3	0	1	2	0
6	H	3	0	1	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	F	4	0	3	0	0
8	A	161	0	0	5	0
8	B	190	0	0	1	0
8	C	132	0	0	6	0
8	D	197	0	0	9	0
8	E	151	0	0	6	0
8	F	161	0	0	4	0
8	G	165	0	0	3	0
8	H	173	0	0	3	0
8	I	181	0	0	2	0
8	J	156	0	0	6	0
8	K	171	0	0	3	0
8	L	171	0	0	3	0
All	All	25215	0	23063	421	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.

The worst 5 of 421 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:34[B]:ARG:HG2	1:A:34[B]:ARG:HH11	1.08	1.16
1:E:52[A]:ARG:HH11	1:E:52[A]:ARG:HG3	1.05	1.14
1:G:244:ARG:HH11	1:G:244:ARG:HG2	1.14	1.11
1:I:144:MSE:SE	1:I:220[B]:VAL:HG23	2.05	1.07
1:D:9:THR:HG21	1:D:14:GLY:HA2	1.33	1.05

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	257/261 (98%)	253 (98%)	3 (1%)	1 (0%)	39	33
1	B	250/261 (96%)	245 (98%)	4 (2%)	1 (0%)	39	33
1	C	252/261 (97%)	247 (98%)	4 (2%)	1 (0%)	39	33
1	D	249/261 (95%)	244 (98%)	4 (2%)	1 (0%)	39	33
1	E	245/261 (94%)	240 (98%)	4 (2%)	1 (0%)	39	33
1	F	255/261 (98%)	247 (97%)	7 (3%)	1 (0%)	39	33
1	G	254/261 (97%)	249 (98%)	3 (1%)	2 (1%)	24	15
1	H	262/261 (100%)	256 (98%)	5 (2%)	1 (0%)	39	33
1	I	257/261 (98%)	251 (98%)	5 (2%)	1 (0%)	39	33
1	J	241/261 (92%)	236 (98%)	4 (2%)	1 (0%)	39	33
1	K	258/261 (99%)	253 (98%)	4 (2%)	1 (0%)	39	33
1	L	247/261 (95%)	240 (97%)	6 (2%)	1 (0%)	39	33
All	All	3027/3132 (97%)	2961 (98%)	53 (2%)	13 (0%)	39	33

5 of 13 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	I	167	TYR
1	J	167	TYR
1	L	167	TYR
1	A	167	TYR
1	B	167	TYR

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	209/198 (106%)	204 (98%)	5 (2%)	57	58
1	B	204/198 (103%)	199 (98%)	5 (2%)	55	55
1	C	206/198 (104%)	193 (94%)	13 (6%)	22	16
1	D	203/198 (102%)	195 (96%)	8 (4%)	39	35

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	199/198 (100%)	188 (94%)	11 (6%)	27	21
1	F	207/198 (104%)	198 (96%)	9 (4%)	35	30
1	G	206/198 (104%)	194 (94%)	12 (6%)	25	19
1	H	214/198 (108%)	204 (95%)	10 (5%)	32	27
1	I	209/198 (106%)	203 (97%)	6 (3%)	50	49
1	J	196/198 (99%)	192 (98%)	4 (2%)	63	65
1	K	210/198 (106%)	201 (96%)	9 (4%)	35	30
1	L	201/198 (102%)	192 (96%)	9 (4%)	34	29
All	All	2464/2376 (104%)	2363 (96%)	101 (4%)	39	32

5 of 101 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	21	ASN
1	G	198	ASN
1	L	65	SER
1	F	198	ASN
1	G	8	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 16 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	198	ASN
1	E	198	ASN
1	I	198	ASN
1	C	198	ASN
1	J	198	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 45 ligands modelled in this entry, 4 are monoatomic - leaving 41 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	GOL	A	260	-	5,5,5	0.35	0	5,5,5	0.28	0
2	GOL	A	270	-	5,5,5	0.40	0	5,5,5	0.25	0
2	GOL	B	261	-	5,5,5	0.34	0	5,5,5	0.31	0
2	GOL	B	270	-	5,5,5	0.44	0	5,5,5	0.27	0
3	URA	B	271	-	4,8,8	1.17	0	6,10,10	9.12	4 (66%)
5	SO4	C	259	-	4,4,4	0.12	0	6,6,6	0.06	0
2	GOL	C	261	-	5,5,5	0.27	0	5,5,5	0.20	0
2	GOL	C	271	-	5,5,5	0.42	0	5,5,5	0.34	0
2	GOL	C	275	-	5,5,5	0.31	0	5,5,5	0.35	0
3	URA	C	276	-	4,8,8	0.49	0	6,10,10	8.14	4 (66%)
2	GOL	D	260	-	5,5,5	0.29	0	5,5,5	0.37	0
3	URA	D	270	-	4,8,8	0.48	0	6,10,10	8.15	4 (66%)
2	GOL	D	271	-	5,5,5	0.43	0	5,5,5	0.93	0
2	GOL	E	259	-	5,5,5	0.45	0	5,5,5	0.72	0
3	URA	E	271	-	4,8,8	0.33	0	6,10,10	7.93	4 (66%)
2	GOL	E	272	-	5,5,5	0.44	0	5,5,5	0.58	0
6	FMT	E	273	-	0,2,2	0.00	-	0,1,1	0.00	-
2	GOL	F	270	-	5,5,5	0.40	0	5,5,5	0.56	0
2	GOL	F	273	-	5,5,5	0.35	0	5,5,5	0.30	0
3	URA	F	274	-	4,8,8	0.59	0	6,10,10	7.32	4 (66%)
7	ACY	F	275	-	1,3,3	2.14	1 (100%)	0,3,3	0.00	-
5	SO4	G	270	-	4,4,4	0.58	0	6,6,6	0.42	0
2	GOL	G	271	-	5,5,5	0.29	0	5,5,5	0.54	0
2	GOL	G	272	-	5,5,5	0.29	0	5,5,5	0.35	0
3	URA	G	273	-	4,8,8	0.73	0	6,10,10	7.84	4 (66%)
2	GOL	G	274	-	5,5,5	0.39	0	5,5,5	0.77	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	URA	H	270	-	4,8,8	0.68	0	6,10,10	7.99	4 (66%)
2	GOL	H	271	-	5,5,5	0.35	0	5,5,5	0.64	0
6	FMT	H	272	-	0,2,2	0.00	-	0,1,1	0.00	-
3	URA	I	259	-	4,8,8	0.62	0	6,10,10	8.04	4 (66%)
2	GOL	I	270	-	5,5,5	0.32	0	5,5,5	0.17	0
5	SO4	I	271	-	4,4,4	0.33	0	6,6,6	0.15	0
2	GOL	I	272	-	5,5,5	0.24	0	5,5,5	0.16	0
2	GOL	J	270	-	5,5,5	0.40	0	5,5,5	0.23	0
3	URA	J	271	-	4,8,8	0.66	0	6,10,10	7.62	4 (66%)
2	GOL	J	273	-	5,5,5	0.31	0	5,5,5	0.34	0
2	GOL	K	270	-	5,5,5	0.40	0	5,5,5	0.23	0
2	GOL	K	271	-	5,5,5	0.43	0	5,5,5	0.65	0
3	URA	K	272	-	4,8,8	0.51	0	6,10,10	7.78	4 (66%)
2	GOL	L	270	-	5,5,5	0.38	0	5,5,5	0.38	0
3	URA	L	271	-	4,8,8	0.59	0	6,10,10	8.18	4 (66%)

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	GOL	A	260	-	-	0/4/4/4	0/0/0/0
2	GOL	A	270	-	-	0/4/4/4	0/0/0/0
2	GOL	B	261	-	-	0/4/4/4	0/0/0/0
2	GOL	B	270	-	-	0/4/4/4	0/0/0/0
3	URA	B	271	-	-	0/0/0/0	0/1/1/1
5	SO4	C	259	-	-	0/0/0/0	0/0/0/0
2	GOL	C	261	-	-	0/4/4/4	0/0/0/0
2	GOL	C	271	-	-	0/4/4/4	0/0/0/0
2	GOL	C	275	-	-	0/4/4/4	0/0/0/0
3	URA	C	276	-	-	0/0/0/0	0/1/1/1
2	GOL	D	260	-	-	0/4/4/4	0/0/0/0
3	URA	D	270	-	-	0/0/0/0	0/1/1/1
2	GOL	D	271	-	-	0/4/4/4	0/0/0/0
2	GOL	E	259	-	-	0/4/4/4	0/0/0/0
3	URA	E	271	-	-	0/0/0/0	0/1/1/1
2	GOL	E	272	-	-	0/4/4/4	0/0/0/0
6	FMT	E	273	-	-	0/0/0/0	0/0/0/0
2	GOL	F	270	-	-	0/4/4/4	0/0/0/0
2	GOL	F	273	-	-	0/4/4/4	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	URA	F	274	-	-	0/0/0/0	0/1/1/1
7	ACY	F	275	-	-	0/0/0/0	0/0/0/0
5	SO4	G	270	-	-	0/0/0/0	0/0/0/0
2	GOL	G	271	-	-	0/4/4/4	0/0/0/0
2	GOL	G	272	-	-	0/4/4/4	0/0/0/0
3	URA	G	273	-	-	0/0/0/0	0/1/1/1
2	GOL	G	274	-	-	0/4/4/4	0/0/0/0
3	URA	H	270	-	-	0/0/0/0	0/1/1/1
2	GOL	H	271	-	-	0/4/4/4	0/0/0/0
6	FMT	H	272	-	-	0/0/0/0	0/0/0/0
3	URA	I	259	-	-	0/0/0/0	0/1/1/1
2	GOL	I	270	-	-	0/4/4/4	0/0/0/0
5	SO4	I	271	-	-	0/0/0/0	0/0/0/0
2	GOL	I	272	-	-	0/4/4/4	0/0/0/0
2	GOL	J	270	-	-	0/4/4/4	0/0/0/0
3	URA	J	271	-	-	0/0/0/0	0/1/1/1
2	GOL	J	273	-	-	0/4/4/4	0/0/0/0
2	GOL	K	270	-	-	0/4/4/4	0/0/0/0
2	GOL	K	271	-	-	0/4/4/4	0/0/0/0
3	URA	K	272	-	-	0/0/0/0	0/1/1/1
2	GOL	L	270	-	-	0/4/4/4	0/0/0/0
3	URA	L	271	-	-	0/0/0/0	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	F	275	ACY	CH3-C	2.14	1.51	1.48

The worst 5 of 44 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	271	URA	N1-C2-N3	-13.29	119.85	128.33
3	K	272	URA	N1-C2-N3	-13.05	120.01	128.33
3	C	276	URA	N1-C2-N3	-12.48	120.37	128.33
3	D	270	URA	N1-C2-N3	-12.45	120.39	128.33
3	I	259	URA	N1-C2-N3	-12.35	120.45	128.33

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

22 monomers are involved in 35 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	270	GOL	1	0
3	B	271	URA	2	0
5	C	259	SO4	1	0
2	C	261	GOL	2	0
2	C	275	GOL	7	0
3	C	276	URA	1	0
2	D	260	GOL	1	0
2	D	271	GOL	1	0
3	E	271	URA	1	0
2	E	272	GOL	1	0
6	E	273	FMT	2	0
2	F	270	GOL	2	0
2	F	273	GOL	3	0
3	F	274	URA	2	0
5	G	270	SO4	1	0
2	H	271	GOL	1	0
6	H	272	FMT	3	0
3	I	259	URA	3	0
2	J	273	GOL	1	0
2	K	270	GOL	2	0
2	K	271	GOL	1	0
3	K	272	URA	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	240/261 (91%)	-0.20	4 (1%) 73 73	11, 20, 38, 87	5 (2%)
1	B	234/261 (89%)	-0.28	3 (1%) 79 80	9, 20, 39, 88	1 (0%)
1	C	238/261 (91%)	-0.08	5 (2%) 67 67	12, 27, 56, 102	4 (1%)
1	D	235/261 (90%)	-0.36	2 (0%) 85 86	11, 20, 43, 89	1 (0%)
1	E	235/261 (90%)	0.01	7 (2%) 54 55	11, 26, 55, 101	5 (2%)
1	F	240/261 (91%)	-0.14	8 (3%) 50 51	11, 22, 60, 110	3 (1%)
1	G	240/261 (91%)	-0.05	6 (2%) 61 61	11, 22, 51, 83	5 (2%)
1	H	239/261 (91%)	-0.03	10 (4%) 40 41	11, 21, 45, 82	1 (0%)
1	I	240/261 (91%)	-0.33	3 (1%) 79 80	11, 22, 45, 71	3 (1%)
1	J	231/261 (88%)	-0.13	6 (2%) 59 60	12, 25, 51, 72	3 (1%)
1	K	241/261 (92%)	-0.37	3 (1%) 81 81	10, 20, 42, 71	4 (1%)
1	L	235/261 (90%)	-0.28	2 (0%) 85 86	10, 20, 42, 75	4 (1%)
All	All	2848/3132 (90%)	-0.19	59 (2%) 67 67	9, 22, 49, 110	39 (1%)

The worst 5 of 59 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	8	LYS	6.0
1	G	9	THR	5.8
1	C	236	ALA	5.8
1	F	232	ILE	5.1
1	H	232	ILE	4.6

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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
3	URA	C	276	8/8	0.84	0.27	6.21	15,22,30,31	8
2	GOL	G	271	6/6	0.90	0.19	4.61	39,45,47,53	0
2	GOL	F	270	6/6	0.87	0.18	3.64	38,46,49,54	0
2	GOL	I	272	6/6	0.87	0.18	3.45	28,30,35,35	0
2	GOL	K	271	6/6	0.94	0.20	3.00	22,27,30,30	0
3	URA	B	271	8/8	0.82	0.21	2.51	31,52,55,59	0
2	GOL	D	271	6/6	0.96	0.17	2.13	16,25,27,28	0
2	GOL	J	273	6/6	0.94	0.14	2.00	45,46,48,49	0
2	GOL	E	272	6/6	0.95	0.14	1.92	33,40,41,46	0
2	GOL	A	260	6/6	0.94	0.14	1.91	35,38,44,48	0
3	URA	D	270	8/8	0.89	0.17	1.77	28,45,49,52	0
6	FMT	E	273	3/3	0.95	0.14	1.60	45,45,46,48	0
2	GOL	J	270	6/6	0.87	0.16	1.48	33,35,37,38	0
2	GOL	C	271	6/6	0.95	0.14	1.47	26,29,30,32	0
2	GOL	B	261	6/6	0.92	0.14	1.41	36,42,45,51	0
2	GOL	B	270	6/6	0.96	0.19	1.35	20,21,23,23	0
2	GOL	E	259	6/6	0.95	0.15	1.13	18,20,23,25	0
2	GOL	A	270	6/6	0.93	0.15	1.12	30,31,32,37	0
2	GOL	G	272	6/6	0.91	0.15	0.87	34,39,42,44	0
2	GOL	F	273	6/6	0.89	0.15	0.81	32,39,41,42	0
2	GOL	C	261	6/6	0.91	0.17	0.68	40,42,44,50	0
2	GOL	G	274	6/6	0.97	0.14	0.67	15,16,19,20	0
2	GOL	I	270	6/6	0.93	0.14	0.49	20,24,29,29	0
2	GOL	C	275	6/6	0.93	0.14	0.48	39,48,54,54	0
3	URA	F	274	8/8	0.94	0.12	0.19	18,26,31,34	0
4	CL	E	270	1/1	0.99	0.09	0.13	43,43,43,43	0
2	GOL	K	270	6/6	0.92	0.11	0.10	28,30,35,35	0
2	GOL	D	260	6/6	0.94	0.11	-0.05	34,36,41,44	0
2	GOL	L	270	6/6	0.94	0.10	-0.13	40,42,42,44	0
3	URA	H	270	8/8	0.96	0.09	-0.17	16,20,23,24	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	URA	J	271	8/8	0.92	0.11	-0.34	28,33,39,39	0
3	URA	L	271	8/8	0.93	0.10	-0.35	19,33,34,36	0
2	GOL	H	271	6/6	0.96	0.10	-0.35	31,36,43,44	0
3	URA	K	272	8/8	0.97	0.08	-0.43	11,15,21,23	0
3	URA	G	273	8/8	0.95	0.10	-0.49	19,26,29,32	0
3	URA	I	259	8/8	0.98	0.08	-0.70	15,17,18,22	0
3	URA	E	271	8/8	0.96	0.10	-0.72	35,40,45,46	0
4	CL	C	274	1/1	0.94	0.10	-	59,59,59,59	0
5	SO4	G	270	5/5	0.97	0.10	-	19,19,24,32	0
5	SO4	C	259	5/5	0.96	0.13	-	37,46,49,53	0
4	CL	I	273	1/1	0.92	0.14	-	72,72,72,72	0
6	FMT	H	272	3/3	0.94	0.19	-	59,59,59,60	0
7	ACY	F	275	4/4	0.86	0.20	-	31,34,37,38	0
5	SO4	I	271	5/5	0.97	0.09	-	29,30,34,41	0
4	CL	I	274	1/1	0.93	0.07	-	54,54,54,54	0

6.5 Other polymers [i](#)

There are no such residues in this entry.