



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:45 AM GMT

PDB ID : 3FWP  
Title : X-ray structure of uridine nucleoside phosphorylase from *Salmonella typhimurium* complexed with phosphate and its inhibitor 2,2'-anhydrouridine at 1.86 Å resolution  
Authors : Lashkov, S.A.; Mikhailov, A.M.; Gabdulkhakov, A.G.  
Deposited on : 2009-01-19  
Resolution : 1.86 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](mailto: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.

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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

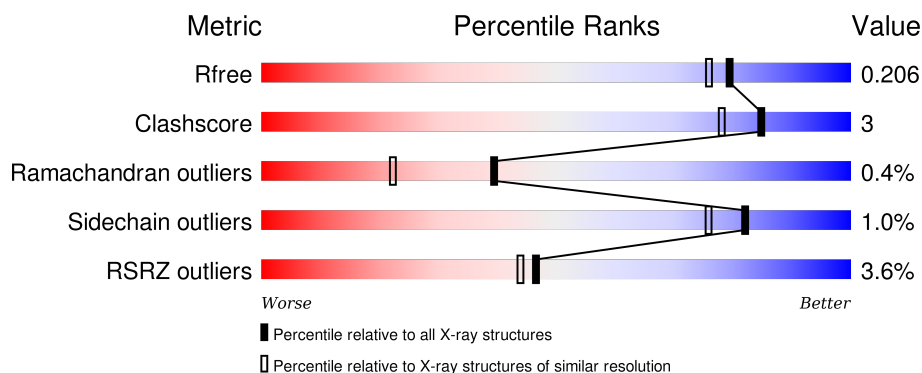
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.86 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	1745 (1.86-1.86)
Clashscore	102246	1898 (1.86-1.86)
Ramachandran outliers	100387	1875 (1.86-1.86)
Sidechain outliers	100360	1875 (1.86-1.86)
RSRZ outliers	91569	1747 (1.86-1.86)

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  $\geq 3$ , 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	253	 3% 91% 7% ••
1	B	253	 2% 88% 8% •
1	C	253	 6% 85% 11% •
1	D	253	 4% 88% 7% 5%
1	E	253	 4% 90% 5% 5%

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Mol	Chain	Length	Quality of chain
1	F	253	

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
4	ANU	B	7016	X	-	-	-
4	ANU	D	7013	X	-	-	-
4	ANU	F	7014	X	-	-	X

## 2 Entry composition [i](#)

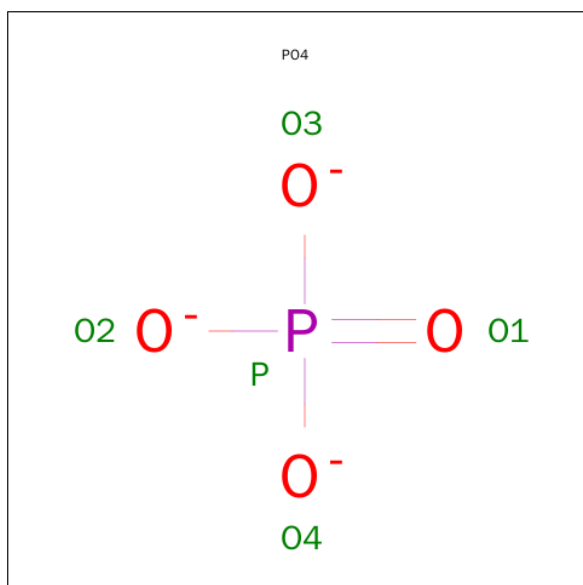
There are 5 unique types of molecules in this entry. The entry contains 12046 atoms, of which 0 are hydrogens and 0 are deuteriums.

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	248	Total	C	N	O	S	0	2	0
			1872	1172	328	360	12			
1	B	242	Total	C	N	O	S	0	2	0
			1822	1143	321	347	11			
1	C	243	Total	C	N	O	S	0	3	0
			1835	1152	323	348	12			
1	D	241	Total	C	N	O	S	0	0	0
			1806	1132	318	344	12			
1	E	240	Total	C	N	O	S	0	5	0
			1823	1146	319	347	11			
1	F	240	Total	C	N	O	S	0	2	0
			1807	1132	321	343	11			

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).

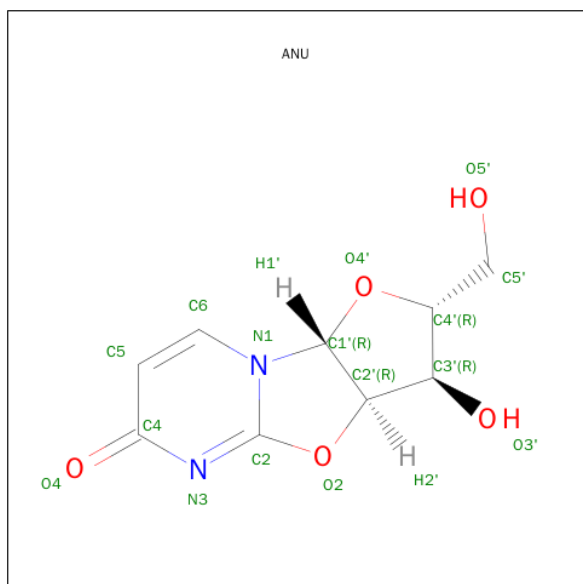


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total O P 5 4 1	0	0
2	D	1	Total O P 5 4 1	0	0
2	F	1	Total O P 5 4 1	0	0

- Molecule 3 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total K 1 1	0	0
3	C	1	Total K 1 1	0	0
3	F	1	Total K 1 1	0	0

- Molecule 4 is 2,2'-ANHYDROURIDINE (three-letter code: ANU) (formula: C<sub>9</sub>H<sub>10</sub>N<sub>2</sub>O<sub>5</sub>).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total C N O 16 9 2 5	0	0
4	D	1	Total C N O 16 9 2 5	0	0
4	F	1	Total C N O 16 9 2 5	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	187	Total 187	O 187	0	0
5	B	187	Total 187	O 187	0	0
5	C	155	Total 155	O 155	0	0
5	D	158	Total 158	O 158	0	0
5	E	179	Total 179	O 179	0	0
5	F	149	Total 149	O 149	0	0

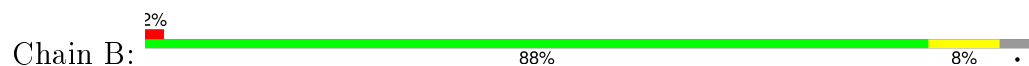
### 3 Residue-property plots [i](#)

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.

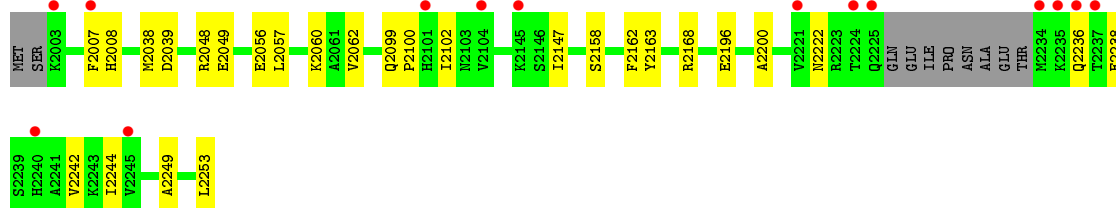
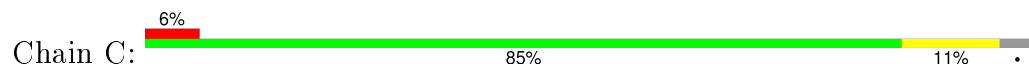
- Molecule 1: Uridine phosphorylase



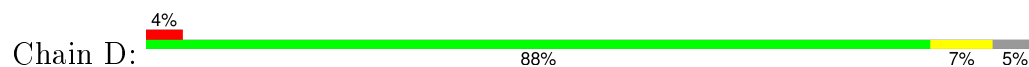
- Molecule 1: Uridine phosphorylase



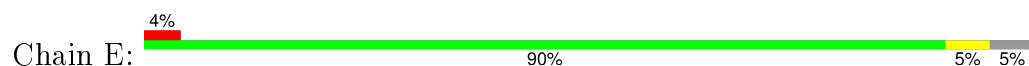
- Molecule 1: Uridine phosphorylase

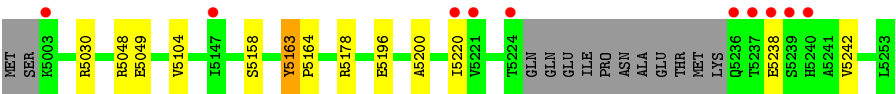


- Molecule 1: Uridine phosphorylase

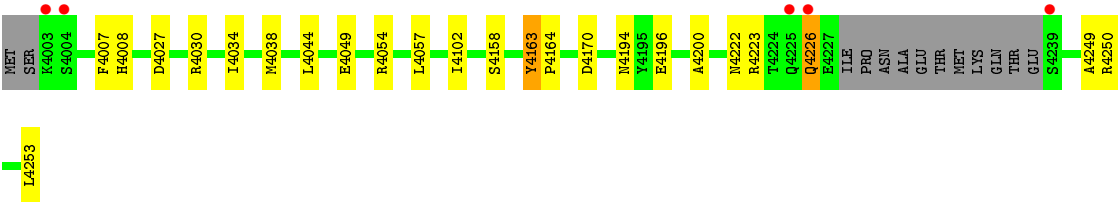
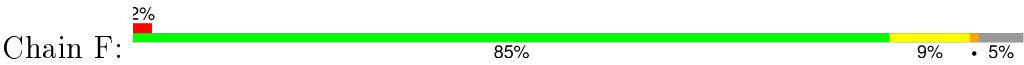


- Molecule 1: Uridine phosphorylase





● Molecule 1: Uridine phosphorylase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	88.79Å 124.07Å 134.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	27.99 – 1.86 27.99 – 1.86	Depositor EDS
% Data completeness (in resolution range)	88.4 (27.99-1.86) 88.4 (27.99-1.86)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.76 (at 1.87Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.176 , 0.206 0.175 , 0.206	Depositor DCC
$R_{free}$ test set	5509 reflections (5.26%)	DCC
Wilson B-factor (Å <sup>2</sup> )	16.9	Xtriage
Anisotropy	0.266	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 53.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 110179 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	12046	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.52% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: ANU, K, PO4

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	0/1906	0.53	0/2579
1	B	0.37	0/1858	0.51	0/2516
1	C	0.36	0/1874	0.51	0/2534
1	D	0.37	0/1834	0.52	0/2482
1	E	0.38	0/1866	0.54	0/2525
1	F	0.36	0/1843	0.52	0/2493
All	All	0.37	0/11181	0.52	0/15129

There are no bond length outliers.

There are no bond angle outliers.

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	1872	0	1886	9	0
1	B	1822	0	1838	9	0
1	C	1835	0	1866	15	0
1	D	1806	0	1821	11	0
1	E	1823	0	1854	8	0
1	F	1807	0	1826	15	0
2	B	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	5	0	0	0	0
2	F	5	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	F	1	0	0	0	0
4	B	16	0	10	0	0
4	D	16	0	10	0	0
4	F	16	0	10	0	0
5	A	187	0	0	0	0
5	B	187	0	0	0	0
5	C	155	0	0	0	0
5	D	158	0	0	1	0
5	E	179	0	0	0	0
5	F	149	0	0	1	0
All	All	12046	0	11121	64	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 3.

All (64) 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:F:4027:ASP:HB3	1:F:4030[B]:ARG:HG3	1.68	0.75
1:B:6125:PRO:HB2	1:B:6127:GLU:OE1	1.95	0.66
1:A:1158:SER:HB3	1:A:1200:ALA:HB2	1.80	0.63
1:C:2158:SER:HB3	1:C:2200:ALA:HB2	1.84	0.59
1:E:5158:SER:HB3	1:E:5200:ALA:HB2	1.84	0.58
1:A:1037:LEU:HD22	1:A:1242:VAL:HG12	1.83	0.58
1:C:2038:MET:HG2	1:C:2057:LEU:HD13	1.85	0.58
1:C:2038:MET:SD	1:C:2062:VAL:HG21	2.44	0.58
1:B:6158:SER:HB3	1:B:6200:ALA:HB2	1.85	0.58
1:B:6108:LEU:HD22	1:B:6152:HIS:HB2	1.86	0.57
1:F:4027:ASP:HB3	1:F:4030[B]:ARG:CG	2.35	0.57
1:F:4158:SER:HB3	1:F:4200:ALA:HB2	1.89	0.53
1:D:3158:SER:HB3	1:D:3200:ALA:HB2	1.90	0.52
1:F:4007:PHE:HD2	1:F:4008:HIS:CE1	2.27	0.52
1:B:6007:PHE:HE1	5:D:943:HOH:O	1.94	0.51
1:A:1038:MET:HG2	1:A:1057:LEU:HD13	1.92	0.51
1:A:1233:THR:HG22	1:A:1235:LYS:H	1.75	0.51
1:B:6049:GLU:HB3	1:D:3049:GLU:HB3	1.94	0.50
1:D:3105:GLY:HA2	1:D:3237:THR:HB	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:4194:ASN:ND2	5:F:1053:HOH:O	2.41	0.49
1:D:3038:MET:HG2	1:D:3057:LEU:HD13	1.93	0.49
1:F:4057:LEU:HG	1:F:4250:ARG:HG2	1.93	0.49
1:E:5030:ARG:NH2	1:E:5238:GLU:HG3	2.28	0.49
1:C:2048:ARG:HB3	1:C:2049:GLU:OE1	2.12	0.49
1:D:3105:GLY:HA2	1:D:3237:THR:CB	2.43	0.48
1:E:5030:ARG:HH21	1:E:5238:GLU:HG3	1.79	0.47
1:C:2162:PHE:O	1:C:2168:ARG:NH1	2.43	0.47
1:C:2007:PHE:HD2	1:C:2008:HIS:CE1	2.32	0.47
1:C:2039:ASP:HB2	1:C:2056:GLU:HB3	1.97	0.47
1:F:4170:ASP:OD1	1:F:4170:ASP:N	2.48	0.47
1:F:4038:MET:HG2	1:F:4057:LEU:HD13	1.97	0.46
1:F:4223:ARG:HA	1:F:4226:GLN:O	2.15	0.46
1:B:6167:GLU:HG2	1:B:6169:TYR:CE1	2.51	0.46
1:A:1049:GLU:HB3	1:F:4049:GLU:HB3	1.97	0.46
1:D:3057:LEU:HB3	1:D:3253:LEU:HD11	1.98	0.45
1:D:3186:GLU:OE2	1:E:5178:ARG:HB2	2.16	0.45
1:D:3238:GLU:O	1:D:3242:VAL:HG23	2.17	0.44
1:E:5163:TYR:HB2	1:E:5164:PRO:CD	2.48	0.44
1:F:4044:LEU:HD11	1:F:4054:ARG:HB2	2.00	0.44
1:A:1232:GLU:HB2	1:A:1233:THR:H	1.59	0.43
1:C:2236:GLN:H	1:C:2236:GLN:CD	2.22	0.43
1:D:3163:TYR:HB2	1:D:3164:PRO:CD	2.49	0.42
1:E:5048:ARG:HB3	1:E:5049:GLU:OE1	2.19	0.42
1:B:6057:LEU:HG	1:B:6250:ARG:HG2	2.01	0.42
1:D:3239:SER:O	1:D:3243:LYS:HG3	2.19	0.42
1:C:2102:ILE:O	1:C:2222:ASN:ND2	2.52	0.42
1:C:2060:LYS:HD2	1:C:2253:LEU:HB3	2.00	0.42
1:E:5238:GLU:O	1:E:5242:VAL:HG23	2.20	0.42
1:C:2147:ILE:HD11	1:C:2244:ILE:HG13	2.02	0.42
1:F:4102:ILE:O	1:F:4222:ASN:ND2	2.48	0.42
1:A:1044:LEU:HD11	1:A:1054:ARG:HB2	2.02	0.41
1:E:5104:VAL:HG13	1:E:5220:ILE:HA	2.02	0.41
1:C:2099:GLN:HA	1:C:2100:PRO:HD3	1.91	0.41
1:C:2238:GLU:O	1:C:2242:VAL:HG23	2.21	0.41
1:A:1163:TYR:HB2	1:A:1164:PRO:HD3	2.03	0.41
1:C:2057:LEU:HB3	1:C:2253:LEU:HD11	2.02	0.41
1:C:2249:ALA:O	1:C:2253:LEU:HG	2.20	0.41
1:D:3049:GLU:HG3	1:D:3068:GLY:HA3	2.03	0.41
1:B:6040:LYS:HE2	1:B:6056:GLU:OE2	2.21	0.41
1:A:1089:PHE:O	1:A:1213:ALA:HA	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:4249:ALA:O	1:F:4253:LEU:HG	2.20	0.40
1:B:6142:GLU:OE1	1:B:6251:ARG:NH2	2.54	0.40
1:F:4030[A]:ARG:HD3	1:F:4034:ILE:HD11	2.02	0.40
1:F:4163:TYR:HB2	1:F:4164:PRO:CD	2.51	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	246/253 (97%)	243 (99%)	2 (1%)	1 (0%)	39	22
1	B	240/253 (95%)	235 (98%)	4 (2%)	1 (0%)	39	22
1	C	242/253 (96%)	236 (98%)	5 (2%)	1 (0%)	39	22
1	D	237/253 (94%)	233 (98%)	3 (1%)	1 (0%)	39	22
1	E	241/253 (95%)	236 (98%)	4 (2%)	1 (0%)	39	22
1	F	238/253 (94%)	235 (99%)	2 (1%)	1 (0%)	39	22
All	All	1444/1518 (95%)	1418 (98%)	20 (1%)	6 (0%)	39	22

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	5163	TYR
1	A	1163	TYR
1	B	6163	TYR
1	D	3163	TYR
1	F	4163	TYR
1	C	2163	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	199/202 (98%)	195 (98%)	4 (2%)	63	47
1	B	194/202 (96%)	192 (99%)	2 (1%)	82	76
1	C	196/202 (97%)	195 (100%)	1 (0%)	92	90
1	D	191/202 (95%)	189 (99%)	2 (1%)	82	76
1	E	195/202 (96%)	194 (100%)	1 (0%)	92	90
1	F	192/202 (95%)	190 (99%)	2 (1%)	82	76
All	All	1167/1212 (96%)	1155 (99%)	12 (1%)	82	76

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1092	ILE
1	A	1196	GLU
1	A	1232	GLU
1	A	1240	HIS
1	B	6145	LYS
1	B	6196	GLU
1	C	2196	GLU
1	D	3092	ILE
1	D	3196	GLU
1	E	5196	GLU
1	F	4196	GLU
1	F	4226	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	6083	GLN
1	B	6209	GLN
1	C	2083	GLN
1	C	2166	GLN
1	D	3166	GLN

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Mol	Chain	Res	Type
1	F	4083	GLN
1	F	4226	GLN

### 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 9 ligands modelled in this entry, 3 are monoatomic - leaving 6 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	PO4	B	7006	-	4,4,4	0.41	0	6,6,6	0.27	0
4	ANU	B	7016	-	13,18,18	3.21	5 (38%)	16,27,27	2.44	5 (31%)
2	PO4	D	7003	-	4,4,4	0.41	0	6,6,6	0.28	0
4	ANU	D	7013	-	13,18,18	3.29	5 (38%)	16,27,27	2.60	7 (43%)
2	PO4	F	7004	-	4,4,4	0.34	0	6,6,6	0.27	0
4	ANU	F	7014	-	13,18,18	3.34	6 (46%)	16,27,27	2.45	7 (43%)

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	PO4	B	7006	-	-	0/0/0/0	0/0/0/0
4	ANU	B	7016	-	1/1/4/4	0/2/26/26	0/2/3/3
2	PO4	D	7003	-	-	0/0/0/0	0/0/0/0
4	ANU	D	7013	-	1/1/4/4	0/2/26/26	0/2/3/3
2	PO4	F	7004	-	-	0/0/0/0	0/0/0/0
4	ANU	F	7014	-	1/1/4/4	0/2/26/26	0/2/3/3

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	7013	ANU	C3'-C2'	-5.65	1.40	1.53
4	F	7014	ANU	C3'-C2'	-5.60	1.40	1.53
4	B	7016	ANU	C3'-C2'	-5.44	1.40	1.53
4	F	7014	ANU	O4'-C4'	-3.96	1.35	1.45
4	B	7016	ANU	O4'-C4'	-3.89	1.36	1.45
4	D	7013	ANU	O4'-C4'	-3.76	1.36	1.45
4	F	7014	ANU	C4-N3	2.02	1.36	1.33
4	B	7016	ANU	O2-C2	2.59	1.39	1.35
4	D	7013	ANU	O2-C2	2.69	1.39	1.35
4	F	7014	ANU	O2-C2	2.84	1.39	1.35
4	B	7016	ANU	C2-N3	4.51	1.41	1.32
4	D	7013	ANU	C2-N3	4.65	1.41	1.32
4	F	7014	ANU	C2-N3	4.93	1.42	1.32
4	B	7016	ANU	O4-C4	7.38	1.42	1.24
4	F	7014	ANU	O4-C4	7.52	1.42	1.24
4	D	7013	ANU	O4-C4	7.55	1.42	1.24

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	7013	ANU	O2-C2'-C3'	-5.67	101.86	110.46
4	B	7016	ANU	O2-C2'-C3'	-4.60	103.48	110.46
4	F	7014	ANU	O2-C2'-C3'	-4.59	103.51	110.46
4	F	7014	ANU	C4'-O4'-C1'	-2.34	107.15	109.72
4	B	7016	ANU	C4'-O4'-C1'	-2.27	107.22	109.72
4	D	7013	ANU	C4'-O4'-C1'	-2.24	107.25	109.72
4	F	7014	ANU	O3'-C3'-C4'	2.05	117.20	111.05
4	D	7013	ANU	C2'-C3'-C4'	2.09	106.79	101.85
4	F	7014	ANU	C2'-C3'-C4'	2.11	106.84	101.85
4	D	7013	ANU	O3'-C3'-C4'	2.24	117.78	111.05
4	B	7016	ANU	O3'-C3'-C2'	2.49	118.34	111.16
4	F	7014	ANU	O3'-C3'-C2'	2.51	118.40	111.16
4	D	7013	ANU	O3'-C3'-C2'	2.72	119.00	111.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	7016	ANU	O4'-C4'-C5'	4.38	118.67	109.17
4	D	7013	ANU	O4'-C4'-C5'	4.52	118.95	109.17
4	F	7014	ANU	O4'-C4'-C5'	4.63	119.20	109.17
4	F	7014	ANU	C4-N3-C2	4.87	118.97	114.14
4	D	7013	ANU	C4-N3-C2	5.33	119.42	114.14
4	B	7016	ANU	C4-N3-C2	5.50	119.59	114.14

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	B	7016	ANU	C2'
4	D	7013	ANU	C2'
4	F	7014	ANU	C2'

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	248/253 (98%)	-0.16	8 (3%)	51	48	8, 15, 34, 61	0
1	B	242/253 (95%)	-0.27	6 (2%)	61	58	9, 15, 28, 33	0
1	C	243/253 (96%)	0.12	14 (5%)	26	25	10, 22, 38, 61	0
1	D	241/253 (95%)	-0.00	10 (4%)	41	38	10, 18, 35, 65	0
1	E	240/253 (94%)	-0.15	10 (4%)	40	37	9, 15, 30, 48	0
1	F	240/253 (94%)	-0.07	5 (2%)	67	65	8, 18, 32, 45	0
All	All	1454/1518 (95%)	-0.09	53 (3%)	46	44	8, 17, 34, 65	0

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	3234	MET	7.3
1	A	1233	THR	6.5
1	A	1232	GLU	6.5
1	A	1234	MET	5.6
1	D	3236	GLN	5.1
1	C	2234	MET	4.7
1	D	3237	THR	4.5
1	F	4226	GLN	4.4
1	A	1228	ILE	4.4
1	D	3104	VAL	4.4
1	D	3235	LYS	4.3
1	D	3105	GLY	4.1
1	E	5236	GLN	3.8
1	C	2236	GLN	3.8
1	C	2104	VAL	3.8
1	C	2003	LYS	3.8
1	E	5237	THR	3.8
1	A	1236	GLN	3.5
1	F	4225	GLN	3.5

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Mol	Chain	Res	Type	RSRZ
1	E	5221	VAL	3.4
1	F	4003	LYS	3.4
1	C	2221	VAL	3.4
1	B	6239	SER	3.3
1	D	3238	GLU	3.3
1	C	2235	LYS	3.3
1	A	1237	THR	3.2
1	C	2101	HIS	3.2
1	C	2007	PHE	3.0
1	E	5238	GLU	2.9
1	F	4004	SER	2.8
1	E	5224	THR	2.8
1	C	2237	THR	2.8
1	E	5220	ILE	2.7
1	C	2225	GLN	2.7
1	C	2240	HIS	2.7
1	E	5003	LYS	2.7
1	D	3101	HIS	2.6
1	B	6148	GLY	2.6
1	C	2145	LYS	2.6
1	B	6004	SER	2.6
1	E	5240	HIS	2.5
1	C	2224	THR	2.5
1	B	6240	HIS	2.3
1	F	4239	SER	2.3
1	D	3148	GLY	2.3
1	E	5147	ILE	2.3
1	A	1235	LYS	2.3
1	E	5239	SER	2.2
1	D	3224	THR	2.2
1	B	6007	PHE	2.2
1	A	1238	GLU	2.1
1	B	6147	ILE	2.0
1	C	2245	VAL	2.0

## 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 [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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
4	ANU	F	7014	16/16	0.82	0.17	2.47	30,31,33,33	0
4	ANU	D	7013	16/16	0.91	0.15	1.25	18,19,19,20	13
3	K	C	8221	1/1	0.95	0.08	-0.21	20,20,20,20	1
4	ANU	B	7016	16/16	0.96	0.08	-0.38	17,17,18,18	0
2	PO4	F	7004	5/5	0.98	0.08	-0.78	27,28,28,28	0
2	PO4	D	7003	5/5	0.97	0.08	-1.04	27,28,28,29	0
2	PO4	B	7006	5/5	0.99	0.06	-2.51	17,17,19,19	0
3	K	B	8001	1/1	0.99	0.04	-3.52	17,17,17,17	1
3	K	F	8002	1/1	0.99	0.03	-4.14	17,17,17,17	1

### 6.5 Other polymers [i](#)

There are no such residues in this entry.