



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:26 PM UTC

PDB ID : 3FWP / pdb_00003fwp
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.

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A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

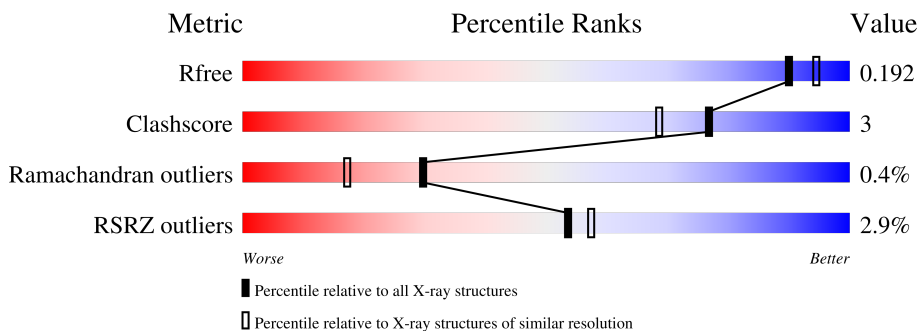
1 Overall quality at a glance

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}	180053	3428 (1.86-1.86)
Clashscore	190562	3579 (1.86-1.86)
Ramachandran outliers	187476	3553 (1.86-1.86)
RSRZ outliers	180081	3429 (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 of 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	253	
1	B	253	
1	C	253	
1	D	253	
1	E	253	
1	F	253	

2 Entry composition [i](#)

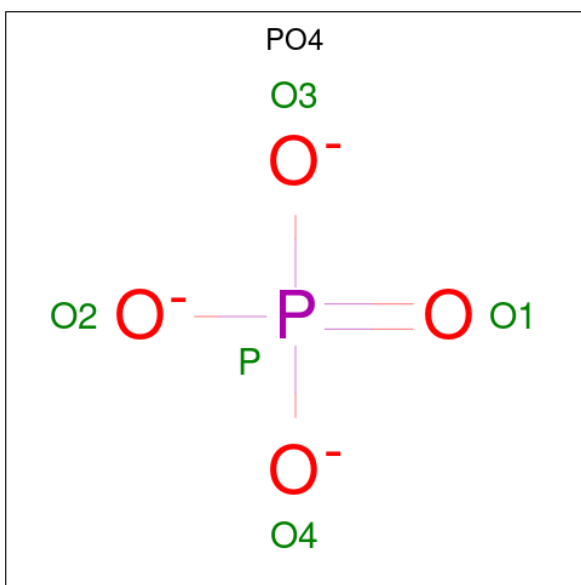
There are 5 unique types of molecules in this entry. The entry contains 12046 atoms, of which 0 are hydrogens and 0 are deuteriums.

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

- Molecule 2 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).

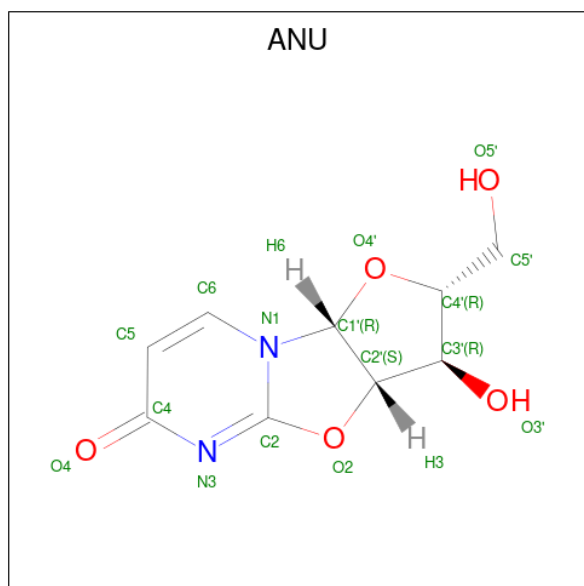


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total O P 5 4 1	0	0
2	D	1	Total O P 5 4 1	0	0
2	F	1	Total O P 5 4 1	0	0

- Molecule 3 is POTASSIUM ION (CCD ID: 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'-Anhydro-(1-beta-D-arabinofuranosyl)uracil (CCD ID: ANU) (formula: C₉H₁₀N₂O₅).



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

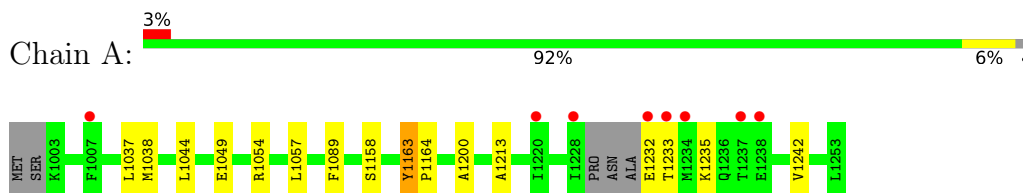
- Molecule 5 is water.

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

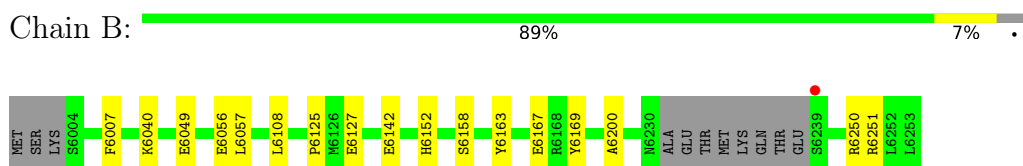
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

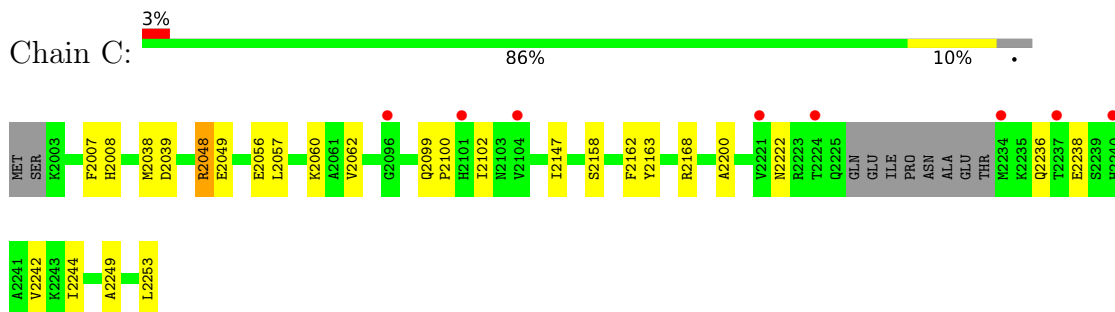
- Molecule 1: Uridine phosphorylase



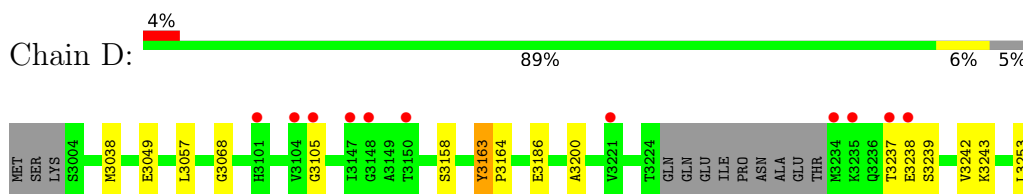
- Molecule 1: Uridine phosphorylase



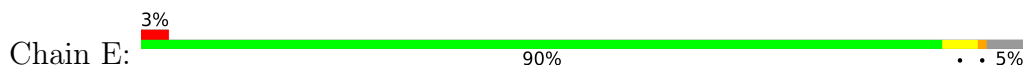
- Molecule 1: Uridine phosphorylase

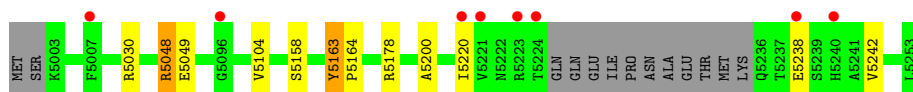


- Molecule 1: Uridine phosphorylase



- Molecule 1: Uridine phosphorylase





- Molecule 1: Uridine phosphorylase

Chain F: 2% 86% 9% 5%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	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
$\langle I/\sigma(I) \rangle$ ¹	2.76 (at 1.87Å)	Xtrriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.176 , 0.206 (Not available) , 0.192	Depositor DCC
R_{free} test set	5509 reflections (4.42%)	wwPDB-VP
Wilson B-factor (Å ²)	16.9	Xtrriage
Anisotropy	0.266	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 45.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12046	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtrriage'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.*

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: PO4, K, ANU

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	0.48	0/1906	0.77	0/2579
1	B	0.48	0/1858	0.75	0/2516
1	C	0.47	0/1874	0.79	1/2534 (0.0%)
1	D	0.48	0/1834	0.78	0/2482
1	E	0.49	0/1866	0.77	1/2525 (0.0%)
1	F	0.47	0/1843	0.77	0/2493
All	All	0.48	0/11181	0.77	2/15129 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	2048	ARG	CB-CA-C	-5.53	110.18	116.54
1	E	5048	ARG	CB-CA-C	-5.12	110.65	116.54

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	10	0
1	C	1835	0	1866	15	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
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
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	65	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 (65) 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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
1:D:3038:MET:HG2	1:D:3057:LEU:HD13	1.93	0.49
1:F:4194:ASN:ND2	5:F:1053:HOH:O	2.41	0.49
1:E:5030:ARG:NH2	1:E:5238:GLU:HG3	2.28	0.49
1:F:4057:LEU:HG	1:F:4250:ARG:HG2	1.93	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:C:2236:GLN:CD	1:C:2236:GLN:H	2.22	0.48
1:C:2162:PHE:O	1:C:2168:ARG:NH1	2.43	0.47
1:E:5030:ARG:HH21	1:E:5238:GLU:HG3	1.79	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:D:3163:TYR:HB2	1:D:3164:PRO:CD	2.49	0.42
1:B:6057:LEU:HG	1:B:6250:ARG:HG2	2.01	0.42
1:E:5048:ARG:HB3	1:E:5049:GLU:OE1	2.19	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:C:2147:ILE:HD11	1:C:2244:ILE:HG13	2.02	0.42
1:E:5238:GLU:O	1:E:5242:VAL:HG23	2.20	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:B:6127:GLU:CD	1:B:6127:GLU:H	2.28	0.41
1:C:2099:GLN:HA	1:C:2100:PRO:HD3	1.91	0.41
1:E:5104:VAL:HG13	1:E:5220:ILE:HA	2.02	0.41
1:A:1163:TYR:HB2	1:A:1164:PRO:HD3	2.03	0.41
1:C:2238:GLU:O	1:C:2242:VAL:HG23	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
1:B:6142:GLU:OE1	1:B:6251:ARG:NH2	2.54	0.40
1:F:4249:ALA:O	1:F:4253:LEU:HG	2.20	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%)	30	17
1	B	240/253 (95%)	235 (98%)	4 (2%)	1 (0%)	30	17
1	C	242/253 (96%)	236 (98%)	5 (2%)	1 (0%)	30	17
1	D	237/253 (94%)	233 (98%)	3 (1%)	1 (0%)	30	17
1	E	241/253 (95%)	236 (98%)	4 (2%)	1 (0%)	30	17
1	F	238/253 (94%)	235 (99%)	2 (1%)	1 (0%)	30	17
All	All	1444/1518 (95%)	1418 (98%)	20 (1%)	6 (0%)	30	17

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	5163	TYR
1	A	1163	TYR
1	B	6163	TYR

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Mol	Chain	Res	Type
1	D	3163	TYR
1	F	4163	TYR
1	C	2163	TYR

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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
4	ANU	F	7014	-	18,18,18	6.07	9 (50%)	23,27,27	3.70	12 (52%)
2	PO4	F	7004	-	4,4,4	0.81	0	6,6,6	0.48	0
2	PO4	D	7003	-	4,4,4	0.90	0	6,6,6	0.56	0
4	ANU	B	7016	-	18,18,18	5.93	10 (55%)	23,27,27	3.61	9 (39%)
2	PO4	B	7006	-	4,4,4	0.89	0	6,6,6	0.52	0
4	ANU	D	7013	-	18,18,18	6.01	10 (55%)	23,27,27	3.71	11 (47%)

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	ANU	F	7014	-	-	2/2/26/26	0/3/3/3
4	ANU	B	7016	-	-	2/2/26/26	0/3/3/3
4	ANU	D	7013	-	-	2/2/26/26	0/3/3/3

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	7013	ANU	C1'-N1	-16.39	1.28	1.46
4	F	7014	ANU	C1'-N1	-16.21	1.28	1.46
4	B	7016	ANU	C1'-N1	-16.12	1.28	1.46
4	F	7014	ANU	C2-N1	14.04	1.56	1.34
4	D	7013	ANU	C2-N1	13.42	1.55	1.34
4	B	7016	ANU	C2-N1	13.33	1.55	1.34
4	D	7013	ANU	O4-C4	7.55	1.42	1.24
4	F	7014	ANU	O4-C4	7.51	1.42	1.24
4	B	7016	ANU	O4-C4	7.37	1.42	1.24
4	F	7014	ANU	C2-N3	6.52	1.42	1.30
4	D	7013	ANU	C2-N3	6.21	1.41	1.30
4	B	7016	ANU	C2-N3	6.06	1.41	1.30
4	D	7013	ANU	C3'-C2'	-5.84	1.40	1.53
4	F	7014	ANU	C3'-C2'	-5.78	1.40	1.53
4	B	7016	ANU	C3'-C2'	-5.62	1.40	1.53
4	F	7014	ANU	C5-C4	5.11	1.55	1.44
4	B	7016	ANU	C5-C4	5.03	1.55	1.44
4	D	7013	ANU	C5-C4	4.94	1.54	1.44
4	F	7014	ANU	O4'-C4'	-4.10	1.35	1.45
4	B	7016	ANU	O4'-C4'	-4.03	1.36	1.45
4	D	7013	ANU	O4'-C4'	-3.90	1.36	1.45
4	F	7014	ANU	O2-C2	3.72	1.39	1.34
4	D	7013	ANU	O2-C2	3.58	1.39	1.34
4	B	7016	ANU	O2-C2	3.50	1.39	1.34
4	B	7016	ANU	C6-C5	2.94	1.41	1.35
4	D	7013	ANU	C6-C5	2.92	1.41	1.35
4	F	7014	ANU	C6-C5	2.92	1.41	1.35
4	D	7013	ANU	O2-C2'	-2.32	1.43	1.46
4	B	7016	ANU	O2-C2'	-2.28	1.43	1.46

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	F	7014	ANU	O2-C2-N3	8.54	133.57	121.24
4	D	7013	ANU	O2-C2-N3	8.35	133.30	121.24
4	B	7016	ANU	O2-C2-N3	8.34	133.28	121.24
4	F	7014	ANU	N1-C2-N3	-8.31	119.83	127.09
4	D	7013	ANU	N1-C2-N3	-8.22	119.91	127.09
4	B	7016	ANU	N1-C2-N3	-8.22	119.92	127.09
4	F	7014	ANU	O2-C2-N1	-7.05	106.59	111.82
4	D	7013	ANU	O2-C2-N1	-6.80	106.78	111.82
4	B	7016	ANU	O2-C2-N1	-6.78	106.79	111.82
4	D	7013	ANU	O2-C2'-C3'	-6.25	101.86	110.62
4	F	7014	ANU	O4'-C1'-N1	-5.74	105.90	111.94
4	D	7013	ANU	O4'-C1'-N1	-5.68	105.96	111.94
4	B	7016	ANU	O4'-C1'-N1	-5.67	105.97	111.94
4	B	7016	ANU	O2-C2'-C3'	-5.09	103.48	110.62
4	F	7014	ANU	O2-C2'-C3'	-5.07	103.51	110.62
4	F	7014	ANU	O4'-C4'-C5'	4.72	119.20	109.22
4	D	7013	ANU	O4'-C4'-C5'	4.61	118.95	109.22
4	B	7016	ANU	O4'-C4'-C5'	4.47	118.67	109.22
4	B	7016	ANU	C2-N3-C4	2.88	119.59	116.66
4	D	7013	ANU	O3'-C3'-C2'	2.79	119.00	111.19
4	D	7013	ANU	C2-N3-C4	2.71	119.42	116.66
4	F	7014	ANU	O3'-C3'-C2'	2.58	118.40	111.19
4	B	7016	ANU	O3'-C3'-C2'	2.56	118.34	111.19
4	D	7013	ANU	O3'-C3'-C4'	2.33	117.78	111.08
4	F	7014	ANU	O4-C4-C5	-2.32	118.68	122.29
4	F	7014	ANU	C2-N3-C4	2.27	118.97	116.66
4	F	7014	ANU	C2'-C3'-C4'	2.25	106.84	101.99
4	F	7014	ANU	C5-C4-N3	2.24	121.26	118.48
4	D	7013	ANU	C2'-C3'-C4'	2.23	106.79	101.99
4	F	7014	ANU	O3'-C3'-C4'	2.13	117.20	111.08
4	D	7013	ANU	C5-C4-N3	2.03	121.00	118.48
4	B	7016	ANU	C5-C4-N3	2.02	120.99	118.48

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	7016	ANU	O4'-C4'-C5'-O5'
4	D	7013	ANU	O4'-C4'-C5'-O5'
4	F	7014	ANU	O4'-C4'-C5'-O5'
4	B	7016	ANU	C3'-C4'-C5'-O5'
4	D	7013	ANU	C3'-C4'-C5'-O5'
4	F	7014	ANU	C3'-C4'-C5'-O5'

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

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	248/253 (98%)	-0.28	8 (3%) 50 54	8, 15, 34, 61	2 (0%)
1	B	242/253 (95%)	-0.27	1 (0%) 88 91	8, 15, 28, 33	2 (0%)
1	C	243/253 (96%)	0.13	8 (3%) 49 53	10, 22, 38, 61	3 (1%)
1	D	241/253 (95%)	0.01	11 (4%) 37 40	10, 18, 35, 65	0
1	E	240/253 (94%)	-0.17	8 (3%) 49 53	7, 15, 30, 48	5 (2%)
1	F	240/253 (94%)	-0.06	6 (2%) 58 62	8, 18, 32, 45	2 (0%)
All	All	1454/1518 (95%)	-0.11	42 (2%) 53 57	7, 17, 34, 65	14 (0%)

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	3105	GLY	4.3
1	E	5220	ILE	4.0
1	E	5221	VAL	3.8
1	D	3237	THR	3.6
1	D	3104	VAL	3.2
1	D	3234	MET	3.0
1	E	5224	THR	3.0
1	A	1232	GLU	3.0
1	C	2101	HIS	3.0
1	F	4226	GLN	2.9
1	D	3101	HIS	2.8
1	E	5007	PHE	2.7
1	F	4007	PHE	2.7
1	C	2104	VAL	2.6
1	D	3238	GLU	2.6
1	D	3235	LYS	2.6
1	A	1233	THR	2.6
1	F	4227	GLU	2.6
1	A	1237	THR	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	1007	PHE	2.5
1	C	2096	GLY	2.5
1	E	5240	HIS	2.5
1	A	1234	MET	2.4
1	C	2221	VAL	2.4
1	C	2240	HIS	2.4
1	F	4004	SER	2.4
1	E	5096	GLY	2.3
1	E	5238	GLU	2.3
1	C	2237	THR	2.3
1	A	1220	ILE	2.3
1	A	1228	ILE	2.3
1	F	4006	VAL	2.3
1	D	3150	THR	2.3
1	F	4003	LYS	2.2
1	B	6239	SER	2.2
1	D	3221	VAL	2.1
1	A	1238	GLU	2.1
1	E	5223	ARG	2.1
1	D	3147	ILE	2.1
1	C	2224	THR	2.1
1	D	3148	GLY	2.1
1	C	2234	MET	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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.

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	ANU	F	7014	16/16	0.83	0.12	30,31,33,33	0
4	ANU	D	7013	16/16	0.91	0.09	18,19,19,20	13
4	ANU	B	7016	16/16	0.95	0.06	17,17,18,18	0
2	PO4	D	7003	5/5	0.95	0.11	27,28,28,29	0
3	K	C	8221	1/1	0.95	0.04	20,20,20,20	1
2	PO4	F	7004	5/5	0.96	0.11	27,28,28,28	0
2	PO4	B	7006	5/5	0.98	0.06	17,17,19,19	0
3	K	F	8002	1/1	0.99	0.05	17,17,17,17	1
3	K	B	8001	1/1	0.99	0.06	17,17,17,17	1

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

There are no such residues in this entry.