



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 3IVD / pdb_00003ivd
Title : Putative 5'-Nucleotidase (c4898) from Escherichia Coli in complex with Uridine
Authors : Ramagopal, U.A.; Toro, R.; Burley, S.K.; Almo, S.C.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2009-08-31
Resolution : 1.90 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 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

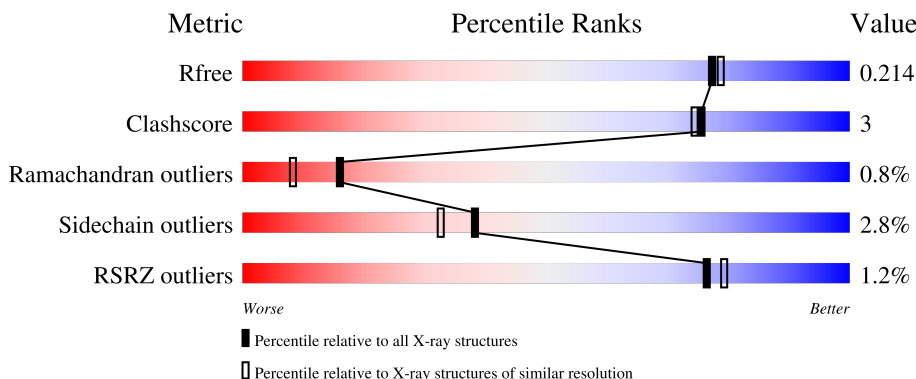
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	509	% 88% 9% ..
1	B	509	% 90% 8% .

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 8223 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 Nucleotidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	499	Total	C	N	O	Se	0	5	0
			3863	2462	651	741	9			
1	B	499	Total	C	N	O	Se	0	2	0
			3860	2455	654	743	8			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	19	MSE	-	expression tag	UNP Q8FBB1
A	20	SER	-	expression tag	UNP Q8FBB1
A	21	LEU	-	expression tag	UNP Q8FBB1
A	520	GLU	-	expression tag	UNP Q8FBB1
A	521	GLY	-	expression tag	UNP Q8FBB1
A	522	HIS	-	expression tag	UNP Q8FBB1
A	523	HIS	-	expression tag	UNP Q8FBB1
A	524	HIS	-	expression tag	UNP Q8FBB1
A	525	HIS	-	expression tag	UNP Q8FBB1
A	526	HIS	-	expression tag	UNP Q8FBB1
A	527	HIS	-	expression tag	UNP Q8FBB1
B	19	MSE	-	expression tag	UNP Q8FBB1
B	20	SER	-	expression tag	UNP Q8FBB1
B	21	LEU	-	expression tag	UNP Q8FBB1
B	520	GLU	-	expression tag	UNP Q8FBB1
B	521	GLY	-	expression tag	UNP Q8FBB1
B	522	HIS	-	expression tag	UNP Q8FBB1
B	523	HIS	-	expression tag	UNP Q8FBB1
B	524	HIS	-	expression tag	UNP Q8FBB1
B	525	HIS	-	expression tag	UNP Q8FBB1
B	526	HIS	-	expression tag	UNP Q8FBB1
B	527	HIS	-	expression tag	UNP Q8FBB1

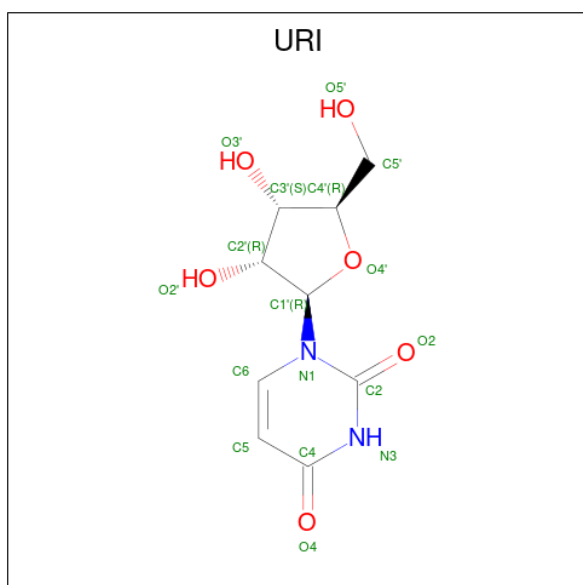
- Molecule 2 is FE (III) ION (CCD ID: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Fe	0	0
			1	1		
2	B	1	Total	Fe	0	0
			1	1		

- Molecule 3 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mn	0	0
			1	1		
3	B	1	Total	Mn	0	0
			1	1		

- Molecule 4 is URIDINE (CCD ID: URI) (formula: C₉H₁₂N₂O₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			17	9	2	6		
4	B	1	Total	C	N	O	0	0
			17	9	2	6		

- Molecule 5 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

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

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	Cl	0	0
			1	1		

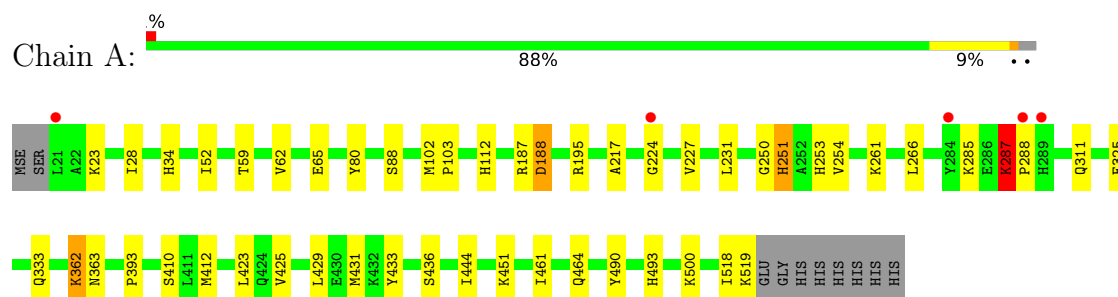
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	230	Total	O	0	0
			230	230		
6	B	230	Total	O	0	0
			230	230		

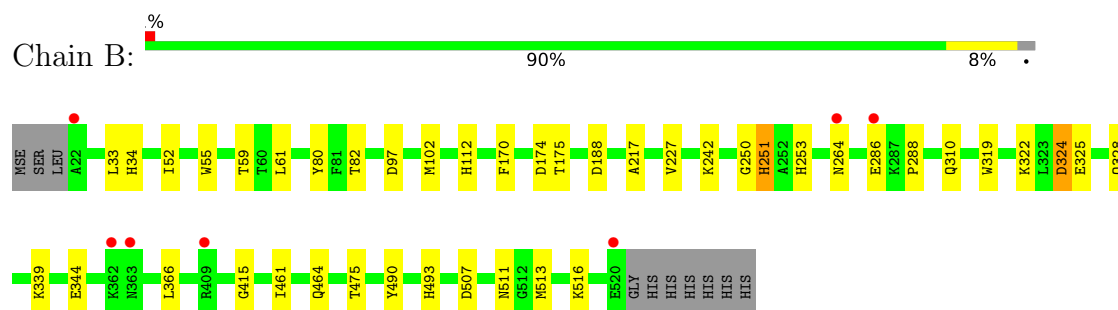
3 Residue-property plots [i](#)

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: Nucleotidase



• Molecule 1: Nucleotidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	47.60Å 67.28Å 86.42Å 74.04° 77.55° 83.76°	Depositor
Resolution (Å)	50.00 – 1.90 50.00 – 1.90	Depositor EDS
% Data completeness (in resolution range)	95.9 (50.00-1.90) 96.1 (50.00-1.90)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.10 (at 1.89Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.167 , 0.210 0.172 , 0.214	Depositor DCC
R_{free} test set	3828 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	21.1	Xtriage
Anisotropy	0.018	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 33.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8223	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.57% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MN, CL, URI, FE

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	0.71	0/3948	0.85	4/5341 (0.1%)
1	B	0.74	0/3936	0.87	1/5324 (0.0%)
All	All	0.72	0/7884	0.86	5/10665 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	287	LYS	CA-C-N	-6.57	111.63	119.84
1	A	287	LYS	C-N-CA	-6.57	111.63	119.84
1	B	475	THR	N-CA-C	6.01	118.62	111.71
1	A	362	LYS	CB-CA-C	-5.76	109.92	116.54
1	A	224	GLY	N-CA-C	5.45	120.97	113.99

There are no chirality outliers.

There are no planarity outliers.

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	3863	0	3861	29	0
1	B	3860	0	3846	18	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	17	0	12	0	0
4	B	17	0	12	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	A	230	0	0	3	0
6	B	230	0	0	1	0
All	All	8223	0	7731	47	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 (47) 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:B:490:TYR:H	1:B:493:HIS:HD2	1.10	0.92
1:A:490:TYR:H	1:A:493:HIS:HD2	1.18	0.86
1:B:490:TYR:H	1:B:493:HIS:CD2	2.01	0.70
1:A:311:GLN:HG2	6:A:685:HOH:O	1.91	0.70
1:A:412:MSE:CE	1:A:423:LEU:HD12	2.26	0.65
1:A:490:TYR:H	1:A:493:HIS:CD2	2.09	0.62
1:A:261:LYS:HG2	1:A:266:LEU:CD2	2.31	0.61
1:A:287:LYS:HB2	6:A:663:HOH:O	2.02	0.59
1:B:59:THR:OG1	1:B:102:MSE:HG2	2.06	0.56
1:A:412:MSE:HE2	1:A:423:LEU:HD12	1.87	0.56
1:A:261:LYS:HG2	1:A:266:LEU:HD23	1.90	0.53
1:B:507:ASP:OD1	1:B:511:ASN:ND2	2.42	0.53
1:A:412:MSE:HE1	1:A:423:LEU:HD12	1.89	0.53
1:A:34:HIS:HE1	1:A:112:HIS:CD2	2.27	0.52
1:B:250:GLY:O	1:B:251:HIS:CB	2.59	0.49
1:A:287:LYS:CB	1:A:288:PRO:HD3	2.43	0.49
1:A:250:GLY:O	1:A:251:HIS:CB	2.60	0.49
1:B:97[A]:ASP:OD2	1:B:319:TRP:HZ2	1.96	0.49
1:B:52:ILE:C	1:B:52:ILE:HD12	2.38	0.48
1:B:344:GLU:HB2	1:B:513:MSE:HE1	1.95	0.48
1:B:61:LEU:C	1:B:61:LEU:HD23	2.38	0.48
1:B:34:HIS:HE1	1:B:112:HIS:CD2	2.32	0.47
1:B:415:GLY:HA3	6:B:565:HOH:O	2.15	0.47
1:A:62:VAL:CG1	1:A:103:PRO:HG2	2.45	0.47
1:A:425:VAL:HB	1:A:429:LEU:HD23	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:62:VAL:HG11	1:A:103:PRO:HG2	1.98	0.45
1:A:52:ILE:HD12	1:A:52:ILE:C	2.43	0.44
1:B:33:LEU:HB3	1:B:55:TRP:CZ2	2.53	0.44
1:A:436:SER:HB3	1:A:519:LYS:HD2	1.99	0.43
1:B:217:ALA:HB2	1:B:253:HIS:CG	2.54	0.43
1:B:324:ASP:O	1:B:328:GLN:HG2	2.18	0.43
1:B:174:ASP:HB3	1:B:175:THR:HG23	2.00	0.43
1:A:362:LYS:HB3	1:A:363:ASN:H	1.62	0.43
1:B:286:GLU:O	1:B:288:PRO:HD3	2.19	0.43
1:A:217:ALA:HB2	1:A:253:HIS:CG	2.53	0.43
1:A:431:MSE:SE	1:A:433:TYR:HB3	2.70	0.42
1:A:444:ILE:HG13	1:A:518:ILE:HD11	2.02	0.42
1:A:217:ALA:HB2	1:A:253:HIS:CD2	2.55	0.42
1:A:88:SER:OG	1:A:393:PRO:O	2.36	0.42
1:B:170:PHE:O	1:B:174:ASP:HB2	2.20	0.42
1:B:97[A]:ASP:OD2	1:B:319:TRP:CZ2	2.73	0.41
1:A:287:LYS:HB2	1:A:288:PRO:HD3	2.02	0.41
1:A:187:ARG:O	1:A:188:ASP:C	2.64	0.41
1:A:28:ILE:HD11	1:A:65:GLU:HG3	2.02	0.41
1:A:231:LEU:HD21	1:A:254:VAL:HB	2.02	0.40
1:A:311:GLN:CG	6:A:685:HOH:O	2.62	0.40
1:A:59:THR:OG1	1:A:102:MSE:HG2	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	502/509 (99%)	488 (97%)	10 (2%)	4 (1%)	16	8
1	B	499/509 (98%)	487 (98%)	8 (2%)	4 (1%)	16	8
All	All	1001/1018 (98%)	975 (97%)	18 (2%)	8 (1%)	16	8

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	188	ASP
1	A	251	HIS
1	B	188	ASP
1	B	251	HIS
1	A	227	VAL
1	A	287	LYS
1	B	82	THR
1	B	227	VAL

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	413/411 (100%)	402 (97%)	11 (3%)	39	34
1	B	412/411 (100%)	399 (97%)	13 (3%)	34	27
All	All	825/822 (100%)	801 (97%)	24 (3%)	38	31

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	LYS
1	A	80	TYR
1	A	195	ARG
1	A	285	LYS
1	A	325	GLU
1	A	333	GLN
1	A	410	SER
1	A	451	LYS
1	A	461	ILE
1	A	464	GLN
1	A	500	LYS
1	B	80	TYR
1	B	242	LYS
1	B	264	ASN
1	B	310[A]	GLN

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Mol	Chain	Res	Type
1	B	310[B]	GLN
1	B	322	LYS
1	B	324	ASP
1	B	325	GLU
1	B	339	LYS
1	B	366	LEU
1	B	461	ILE
1	B	464	GLN
1	B	516	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	120	ASN
1	A	124	GLN
1	A	240	GLN
1	A	290	ASN
1	A	295	ASN
1	A	314	GLN
1	A	320	ASN
1	A	449	ASN
1	A	493	HIS
1	B	120	ASN
1	B	124	GLN
1	B	236	GLN
1	B	240	GLN
1	B	264	ASN
1	B	290	ASN
1	B	295	ASN
1	B	320	ASN
1	B	329	GLN
1	B	493	HIS

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 6 are monoatomic - leaving 2 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	URI	B	603	-	18,18,18	0.53	0	26,26,26	0.86	0
4	URI	A	603	-	18,18,18	0.71	1 (5%)	26,26,26	1.11	1 (3%)

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	URI	B	603	-	-	0/6/22/22	0/2/2/2
4	URI	A	603	-	-	0/6/22/22	0/2/2/2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	603	URI	O3'-C3'	2.00	1.47	1.43

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	603	URI	C4-N3-C2	-3.06	122.82	126.61

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	491/509 (96%)	-0.25	5 (1%) 79 82	11, 20, 35, 50	4 (0%)
1	B	491/509 (96%)	-0.29	7 (1%) 73 76	10, 20, 35, 52	3 (0%)
All	All	982/1018 (96%)	-0.27	12 (1%) 76 79	10, 20, 35, 52	7 (0%)

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	21	LEU	4.0
1	B	22	ALA	3.5
1	A	289	HIS	3.1
1	B	264	ASN	3.1
1	A	288	PRO	2.7
1	B	286	GLU	2.6
1	B	362	LYS	2.3
1	A	284	TYR	2.2
1	B	409	ARG	2.2
1	B	363	ASN	2.1
1	B	520	GLU	2.1
1	A	224	GLY	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.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	URI	A	603	17/17	0.96	0.05	14,17,19,19	0
4	URI	B	603	17/17	0.97	0.04	14,17,19,19	0
3	MN	A	602	1/1	0.99	0.04	23,23,23,23	0
5	CL	A	604	1/1	0.99	0.06	19,19,19,19	0
5	CL	B	604	1/1	0.99	0.03	16,16,16,16	0
2	FE	A	601	1/1	1.00	0.02	19,19,19,19	0
3	MN	B	602	1/1	1.00	0.02	19,19,19,19	0
2	FE	B	601	1/1	1.00	0.01	15,15,15,15	0

6.5 Other polymers [i](#)

There are no such residues in this entry.