



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 12:16 PM UTC

PDB ID : 4TXL / pdb_00004txl
Title : Crystal structure of uridine phosphorylase from *Schistosoma mansoni* in complex with uracil
Authors : Marinho, A.; Torini, J.; Romanello, L.; Cassago, A.; DeMarco, R.; Brandao-Neto, J.; Pereira, H.M.
Deposited on : 2014-07-03
Resolution : 1.92 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

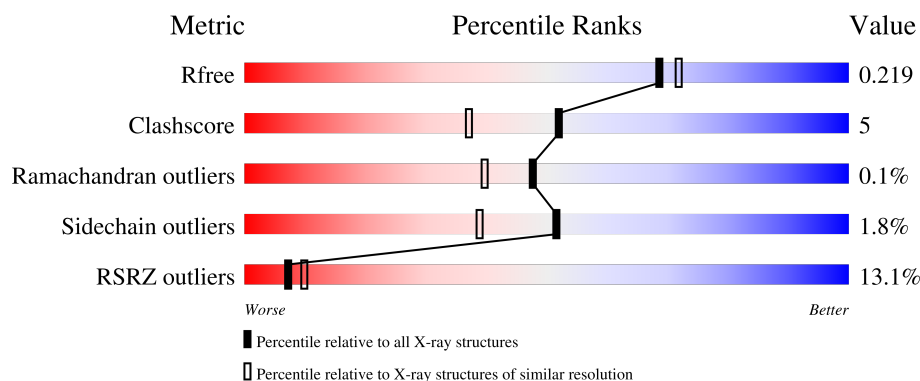
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.92 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	1188 (1.92-1.92)
Clashscore	190562	1209 (1.92-1.92)
Ramachandran outliers	187476	1195 (1.92-1.92)
Sidechain outliers	187428	1195 (1.92-1.92)
RSRZ outliers	180081	1188 (1.92-1.92)

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	296	10% 87% 8% 5%
1	B	296	39% 76% 19% • •
1	C	296	% 93% 5% •
1	D	296	% 95% • •

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
3	URA	A	302	-	X	-	-
3	URA	B	302	-	X	-	-
3	URA	C	302	-	X	-	-
3	URA	D	302	-	X	-	-

2 Entry composition [i](#)

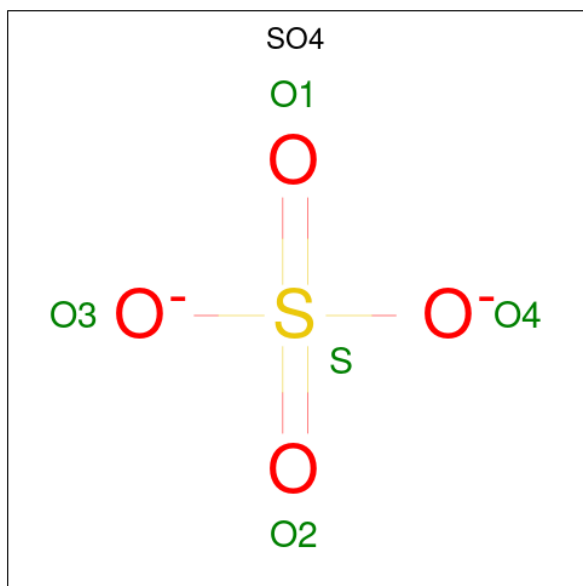
There are 4 unique types of molecules in this entry. The entry contains 9748 atoms, of which 0 are hydrogens and 0 are deuteriums.

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	C	290	Total	C	N	O	S	0	1	0
			2217	1409	383	411	14			
1	D	289	Total	C	N	O	S	0	1	0
			2203	1398	384	407	14			
1	A	282	Total	C	N	O	S	0	1	0
			2083	1324	359	386	14			
1	B	284	Total	C	N	O	S	0	0	0
			2054	1297	357	387	13			

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



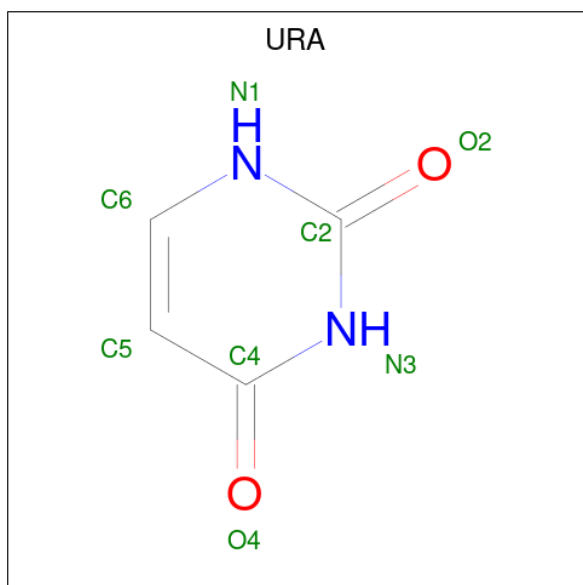
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is URACIL (CCD ID: URA) (formula: C₄H₄N₂O₂).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	365	Total	O	0	0
			365	365		
4	D	341	Total	O	0	0
			341	341		
4	A	215	Total	O	0	0
			215	215		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	218	Total 218	O 218	0	0

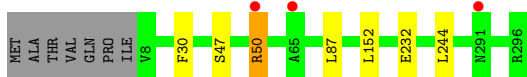
3 Residue-property plots [i](#)

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.

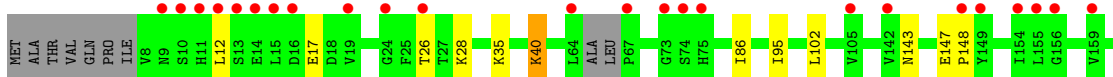
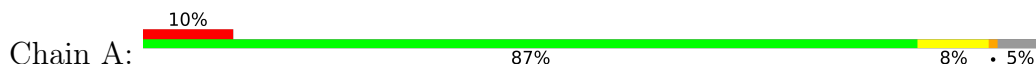
- Molecule 1: Uridine phosphorylase



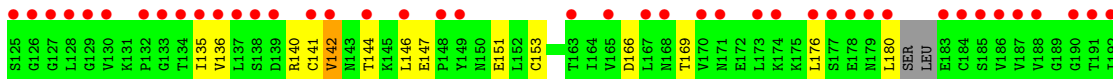
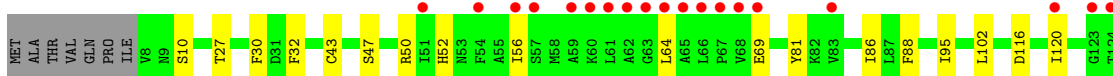
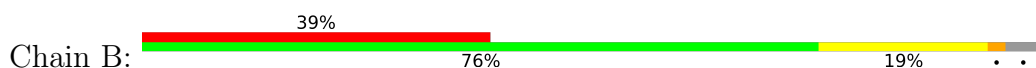
- Molecule 1: Uridine phosphorylase

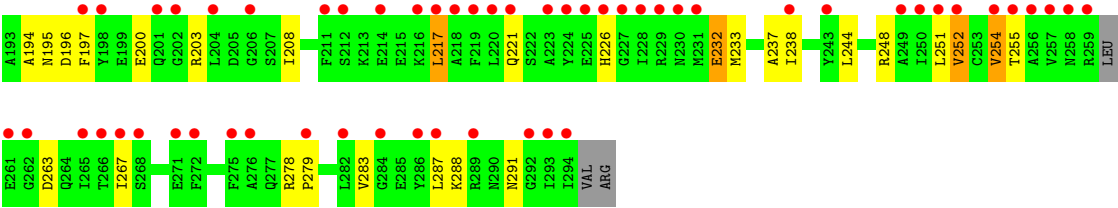


- Molecule 1: Uridine phosphorylase



- Molecule 1: Uridine phosphorylase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	95.85Å 108.53Å 118.67Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	71.84 – 1.92 71.84 – 1.92	Depositor EDS
% Data completeness (in resolution range)	99.6 (71.84-1.92) 99.6 (71.84-1.92)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.24 (at 1.92Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE: 1.9_1692)	Depositor
R, R_{free}	0.191 , 0.220 0.193 , 0.219	Depositor DCC
R_{free} test set	4744 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	23.7	Xtriage
Anisotropy	0.054	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 58.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9748	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.40% 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: URA, SO4

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	0.32	0/2122	0.70	0/2874
1	B	0.35	0/2092	0.77	0/2839
1	C	0.37	0/2260	0.69	0/3058
1	D	0.37	0/2243	0.69	0/3035
All	All	0.35	0/8717	0.71	0/11806

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	2083	0	1989	19	0
1	B	2054	0	1882	52	0
1	C	2217	0	2183	12	0
1	D	2203	0	2161	3	0
2	A	5	0	0	0	0
2	B	5	0	0	1	0
2	C	5	0	0	0	0
2	D	5	0	0	0	0
3	A	8	0	3	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	8	0	3	0	0
3	C	8	0	3	0	0
3	D	8	0	3	0	0
4	A	215	0	0	10	5
4	B	218	0	0	30	0
4	C	365	0	0	7	2
4	D	341	0	0	0	3
All	All	9748	0	8227	85	5

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 5.

All (85) 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:140:ARG:NH2	4:B:403:HOH:O	2.06	0.88
1:A:35:LYS:O	4:A:543:HOH:O	1.92	0.88
1:B:196:ASP:OD1	4:B:575:HOH:O	1.91	0.88
4:A:597:HOH:O	1:B:233:MET:SD	2.36	0.84
1:B:141:CYS:SG	4:B:428:HOH:O	2.35	0.82
1:B:153:CYS:O	4:B:423:HOH:O	1.99	0.81
1:A:12:LEU:O	4:A:582:HOH:O	2.00	0.80
1:C:171:ASN:OD1	4:C:729:HOH:O	2.01	0.77
1:A:26:THR:HG23	1:A:28:LYS:H	1.48	0.76
2:B:301:SO4:S	4:B:573:HOH:O	2.44	0.75
1:B:64:LEU:O	4:B:581:HOH:O	2.04	0.75
1:C:151:GLU:OE2	4:C:687:HOH:O	2.08	0.70
1:B:200:GLU:OE1	4:B:607:HOH:O	2.09	0.70
1:B:255:THR:O	4:B:610:HOH:O	2.10	0.69
1:B:50:ARG:NH1	4:B:573:HOH:O	2.26	0.69
1:B:291:ASN:ND2	4:B:580:HOH:O	2.09	0.69
1:B:140:ARG:NH1	1:B:226:HIS:O	2.26	0.68
1:C:168:ASN:ND2	4:C:742:HOH:O	2.27	0.67
1:B:232:GLU:HB2	4:B:452:HOH:O	1.93	0.67
1:B:151:GLU:O	4:B:609:HOH:O	2.13	0.67
1:B:203:ARG:HG2	1:B:263:ASP:HB3	1.77	0.67
1:A:17:GLU:OE2	4:A:578:HOH:O	2.15	0.65
1:B:254:VAL:HG22	4:B:610:HOH:O	1.97	0.65
1:C:261:GLU:OE2	4:C:714:HOH:O	2.15	0.65
1:A:281:GLN:NE2	4:A:530:HOH:O	2.29	0.65
1:B:116:ASP:OD2	4:B:595:HOH:O	2.14	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:200:GLU:HB2	4:B:607:HOH:O	1.95	0.64
1:A:168:ASN:OD1	1:A:168:ASN:N	2.32	0.63
1:B:196:ASP:O	4:B:587:HOH:O	2.16	0.59
1:B:251:LEU:HB2	4:B:428:HOH:O	2.02	0.58
1:B:288:LYS:O	4:B:547:HOH:O	2.17	0.58
1:B:120:ILE:HD13	1:B:283:VAL:HG11	1.85	0.57
1:A:86:ILE:HD13	1:A:287:LEU:HD13	1.87	0.57
1:C:208:ILE:HD13	1:D:244:LEU:HD23	1.87	0.56
1:C:183:GLU:H	1:C:183:GLU:CD	2.14	0.56
1:C:285:GLU:OE1	4:C:734:HOH:O	2.18	0.55
1:B:217:LEU:HB3	1:B:221:GLN:HE22	1.72	0.55
1:B:200:GLU:CB	4:B:607:HOH:O	2.53	0.55
1:C:183:GLU:HG3	4:C:433:HOH:O	2.06	0.54
1:B:166:ASP:HB2	4:B:612:HOH:O	2.08	0.53
1:A:102:LEU:HD11	1:A:238:ILE:HD13	1.91	0.53
1:B:208:ILE:O	4:B:563:HOH:O	2.18	0.53
1:A:235:GLY:HA3	4:A:461:HOH:O	2.09	0.53
1:B:95:ILE:HG23	1:B:237:ALA:HB2	1.91	0.52
1:A:95:ILE:HG23	1:A:237:ALA:HB2	1.92	0.51
1:B:194:ALA:HB1	4:B:607:HOH:O	2.10	0.51
1:A:243:TYR:N	4:A:476:HOH:O	2.43	0.50
1:B:102:LEU:HD11	1:B:238:ILE:HG12	1.94	0.50
1:B:248:ARG:NH2	4:B:615:HOH:O	2.40	0.50
1:B:169:THR:O	4:B:558:HOH:O	2.19	0.49
1:B:255:THR:N	4:B:610:HOH:O	2.43	0.48
1:A:196:ASP:OD1	1:A:197:PHE:N	2.46	0.48
1:B:10:SER:OG	4:B:401:HOH:O	1.98	0.47
1:B:81:TYR:HB2	1:B:88:PHE:HB2	1.95	0.47
1:B:153:CYS:SG	4:B:556:HOH:O	2.61	0.47
1:A:208:ILE:HD13	1:B:244:LEU:HD23	1.97	0.46
1:B:203:ARG:HG2	1:B:263:ASP:CB	2.46	0.46
1:B:47:SER:HB3	1:B:50:ARG:HB2	1.98	0.45
1:B:195:ASN:HA	4:B:430:HOH:O	2.15	0.45
1:B:195:ASN:N	4:B:607:HOH:O	2.49	0.45
1:A:40:LYS:HD3	4:A:521:HOH:O	2.16	0.45
1:B:27:THR:HA	1:B:32:PHE:CE1	2.52	0.45
4:C:649:HOH:O	1:D:152:LEU:HD23	2.17	0.44
1:B:252:VAL:HG22	1:B:279:PRO:HB3	2.00	0.44
1:B:52:HIS:CD2	1:B:81:TYR:CZ	3.06	0.44
1:B:142:VAL:HG22	1:B:147:GLU:C	2.43	0.44
1:C:39:VAL:HG21	1:C:87:LEU:HG	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:143:ASN:HB3	4:A:495:HOH:O	2.18	0.43
1:B:135:ILE:HD11	1:B:278:ARG:HB3	2.00	0.43
1:B:176:LEU:O	1:B:180:LEU:HG	2.19	0.43
1:B:267:ILE:O	4:B:454:HOH:O	2.21	0.43
1:C:95:ILE:HG23	1:C:237:ALA:HB2	2.01	0.43
1:B:43:CYS:SG	1:B:120:ILE:HD11	2.58	0.43
1:B:196:ASP:OD1	1:B:197:PHE:N	2.52	0.42
1:B:217:LEU:HD13	1:B:217:LEU:HA	1.86	0.42
1:A:289:ARG:HD2	4:A:401:HOH:O	2.18	0.42
1:B:69:GLU:CD	1:B:69:GLU:H	2.28	0.42
1:D:47:SER:HB3	1:D:50[A]:ARG:HB2	2.02	0.41
1:A:196:ASP:HB3	1:A:200:GLU:HB2	2.03	0.41
1:B:142:VAL:HG13	1:B:146:LEU:HA	2.02	0.41
1:C:196:ASP:OD1	1:C:197:PHE:N	2.54	0.41
1:A:147:GLU:HA	1:A:148:PRO:HD3	1.88	0.41
1:C:51:ILE:HG13	1:C:88:PHE:HB3	2.03	0.40
1:A:40:LYS:HD2	1:A:40:LYS:HA	1.83	0.40
1:B:86:ILE:HD13	1:B:287:LEU:HD13	2.03	0.40

All (5) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:440:HOH:O	4:A:403:HOH:O[4_455]	2.02	0.18
4:C:438:HOH:O	4:A:414:HOH:O[3_545]	2.12	0.08
4:C:438:HOH:O	4:A:419:HOH:O[3_545]	2.16	0.04
4:D:414:HOH:O	4:A:418:HOH:O[1_455]	2.16	0.04
4:D:439:HOH:O	4:A:405:HOH:O[1_455]	2.16	0.04

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	277/296 (94%)	270 (98%)	6 (2%)	1 (0%)	30	19
1	B	278/296 (94%)	271 (98%)	7 (2%)	0	100	100
1	C	289/296 (98%)	280 (97%)	9 (3%)	0	100	100
1	D	288/296 (97%)	281 (98%)	7 (2%)	0	100	100
All	All	1132/1184 (96%)	1102 (97%)	29 (3%)	1 (0%)	48	40

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	166	ASP

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	211/252 (84%)	209 (99%)	2 (1%)	70	67
1	B	199/252 (79%)	190 (96%)	9 (4%)	24	9
1	C	236/252 (94%)	235 (100%)	1 (0%)	84	83
1	D	233/252 (92%)	228 (98%)	5 (2%)	47	33
All	All	879/1008 (87%)	862 (98%)	17 (2%)	51	37

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	30	PHE
1	D	30	PHE
1	D	50[A]	ARG
1	D	50[B]	ARG
1	D	87	LEU
1	D	232	GLU
1	A	40	LYS
1	A	168	ASN
1	B	30	PHE
1	B	56	ILE

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Mol	Chain	Res	Type
1	B	136	VAL
1	B	142	VAL
1	B	144	THR
1	B	217	LEU
1	B	232	GLU
1	B	252	VAL
1	B	254	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	11	HIS
1	A	52	HIS
1	A	281	GLN
1	A	290	ASN
1	B	110	HIS
1	B	221	GLN

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](#)

8 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the

expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	URA	D	302	-	8,8,8	4.56	6 (75%)	10,10,10	2.84	5 (50%)
3	URA	A	302	-	8,8,8	4.51	6 (75%)	10,10,10	2.88	5 (50%)
2	SO4	D	301	-	4,4,4	0.25	0	6,6,6	0.10	0
3	URA	C	302	-	8,8,8	4.47	6 (75%)	10,10,10	2.82	6 (60%)
3	URA	B	302	-	8,8,8	4.69	6 (75%)	10,10,10	2.84	5 (50%)
2	SO4	A	301	-	4,4,4	0.25	0	6,6,6	0.24	0
2	SO4	C	301	-	4,4,4	0.23	0	6,6,6	0.12	0
2	SO4	B	301	-	4,4,4	0.25	0	6,6,6	0.18	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	URA	B	302	-	-	-	0/1/1/1
3	URA	A	302	-	-	-	0/1/1/1
3	URA	C	302	-	-	-	0/1/1/1
3	URA	D	302	-	-	-	0/1/1/1

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	302	URA	O4-C4	8.51	1.41	1.24
3	B	302	URA	O2-C2	8.30	1.41	1.23
3	D	302	URA	O2-C2	8.20	1.40	1.23
3	C	302	URA	O2-C2	8.14	1.40	1.23
3	D	302	URA	O4-C4	8.05	1.40	1.24
3	A	302	URA	O2-C2	8.05	1.40	1.23
3	A	302	URA	O4-C4	8.01	1.40	1.24
3	C	302	URA	O4-C4	7.79	1.39	1.24
3	D	302	URA	C2-N1	-3.30	1.32	1.36
3	C	302	URA	C2-N3	-3.17	1.32	1.37
3	A	302	URA	C2-N1	-3.10	1.32	1.36
3	B	302	URA	C2-N3	-3.03	1.32	1.37
3	B	302	URA	C2-N1	-3.02	1.32	1.36
3	D	302	URA	C2-N3	-3.00	1.32	1.37
3	B	302	URA	C6-C5	2.90	1.41	1.35
3	C	302	URA	C2-N1	-2.84	1.33	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	302	URA	C2-N3	-2.80	1.32	1.37
3	A	302	URA	C4-N3	-2.78	1.33	1.38
3	D	302	URA	C6-C5	2.74	1.41	1.35
3	A	302	URA	C6-C5	2.67	1.40	1.35
3	C	302	URA	C6-C5	2.60	1.40	1.35
3	B	302	URA	C4-N3	-2.38	1.34	1.38
3	C	302	URA	C4-N3	-2.38	1.34	1.38
3	D	302	URA	C4-N3	-2.16	1.34	1.38

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	302	URA	N1-C2-N3	5.76	121.25	115.17
3	A	302	URA	N1-C2-N3	5.71	121.19	115.17
3	D	302	URA	N1-C2-N3	5.52	120.99	115.17
3	C	302	URA	N1-C2-N3	5.42	120.89	115.17
3	D	302	URA	C4-N3-C2	-3.87	121.80	125.55
3	A	302	URA	C4-N3-C2	-3.67	121.99	125.55
3	C	302	URA	C4-N3-C2	-3.60	122.06	125.55
3	B	302	URA	C4-N3-C2	-3.49	122.16	125.55
3	A	302	URA	O2-C2-N1	-3.35	119.34	122.79
3	B	302	URA	O2-C2-N1	-3.13	119.56	122.79
3	D	302	URA	O2-C2-N1	-3.09	119.60	122.79
3	C	302	URA	C5-C4-N3	2.95	118.93	114.80
3	D	302	URA	C5-C4-N3	2.86	118.81	114.80
3	C	302	URA	O4-C4-C5	-2.77	120.39	125.16
3	A	302	URA	C5-C4-N3	2.75	118.66	114.80
3	C	302	URA	O2-C2-N1	-2.64	120.06	122.79
3	D	302	URA	O4-C4-C5	-2.63	120.63	125.16
3	B	302	URA	C5-C4-N3	2.59	118.43	114.80
3	A	302	URA	O4-C4-C5	-2.53	120.80	125.16
3	B	302	URA	O4-C4-C5	-2.42	120.99	125.16
3	C	302	URA	C6-C5-C4	-2.24	117.34	119.30

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	301	SO4	1	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	282/296 (95%)	0.65	29 (10%) 12 16	13, 39, 59, 72	1 (0%)
1	B	284/296 (95%)	1.71	116 (40%) 0 0	17, 53, 77, 92	0
1	C	290/296 (97%)	-0.41	2 (0%) 84 88	8, 19, 33, 47	1 (0%)
1	D	289/296 (97%)	-0.28	3 (1%) 79 85	8, 22, 41, 58	1 (0%)
All	All	1145/1184 (96%)	0.41	150 (13%) 7 10	8, 28, 67, 92	3 (0%)

All (150) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	66	LEU	5.4
1	B	180	LEU	5.3
1	B	228	ILE	5.1
1	B	56	ILE	4.9
1	B	130	VAL	4.8
1	B	136	VAL	4.7
1	B	176	LEU	4.7
1	B	128	LEU	4.5
1	B	135	ILE	4.5
1	B	190	GLY	4.4
1	B	186	VAL	4.4
1	D	50[A]	ARG	4.4
1	B	229	ARG	4.3
1	B	146	LEU	4.3
1	B	139	ASP	4.3
1	B	294	ILE	4.1
1	B	191	THR	4.1
1	B	142	VAL	4.1
1	B	224	TYR	4.1
1	B	67	PRO	4.1
1	A	64	LEU	4.1

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Mol	Chain	Res	Type	RSRZ
1	B	220	LEU	4.0
1	B	221	GLN	3.9
1	B	252	VAL	3.9
1	B	204	LEU	3.9
1	B	217	LEU	3.9
1	B	134	THR	3.8
1	B	65	ALA	3.8
1	B	59	ALA	3.7
1	B	254	VAL	3.6
1	B	173	LEU	3.6
1	B	165	VAL	3.5
1	B	249	ALA	3.5
1	C	181	SER	3.5
1	A	9	ASN	3.5
1	B	68	VAL	3.5
1	B	133	GLY	3.4
1	A	244	LEU	3.4
1	A	75	HIS	3.4
1	B	51	ILE	3.4
1	B	132	PRO	3.4
1	B	202	GLY	3.3
1	B	227	GLY	3.3
1	B	286	TYR	3.3
1	B	267	ILE	3.3
1	B	83	VAL	3.3
1	B	279	PRO	3.3
1	B	184	CYS	3.3
1	A	19	VAL	3.2
1	B	214	GLU	3.2
1	B	258	ASN	3.2
1	B	284	GLY	3.2
1	B	250	ILE	3.2
1	B	275	PHE	3.2
1	B	223	ALA	3.1
1	B	149	TYR	3.1
1	B	168	ASN	3.1
1	B	64	LEU	3.1
1	B	211	PHE	3.1
1	B	137	LEU	3.0
1	B	129	GLY	3.0
1	B	225	GLU	3.0
1	B	219	PHE	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	257	VAL	2.9
1	A	156	GLY	2.9
1	B	255	THR	2.9
1	B	272	PHE	2.9
1	B	256	ALA	2.9
1	B	287	LEU	2.9
1	B	126	GLY	2.9
1	B	123	GLY	2.9
1	B	201	GLN	2.8
1	A	15	LEU	2.8
1	B	265	ILE	2.8
1	A	159	VAL	2.8
1	B	183	GLU	2.8
1	A	11	HIS	2.8
1	A	149	TYR	2.8
1	B	69	GLU	2.8
1	A	74	SER	2.7
1	B	120	ILE	2.7
1	B	188	VAL	2.7
1	B	231	MET	2.7
1	B	57	SER	2.7
1	B	54	PHE	2.7
1	A	155	LEU	2.7
1	B	141	CYS	2.6
1	B	271	GLU	2.6
1	C	7	ILE	2.6
1	B	230	ASN	2.6
1	A	154	ILE	2.6
1	B	192	ILE	2.6
1	B	216	LYS	2.6
1	D	65	ALA	2.6
1	B	62	ALA	2.6
1	B	187	VAL	2.6
1	B	218	ALA	2.6
1	B	148	PRO	2.5
1	A	67	PRO	2.5
1	A	182	LEU	2.5
1	B	167	LEU	2.5
1	B	292	GLY	2.5
1	A	246	GLY	2.5
1	A	16	ASP	2.5
1	A	12	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	138	SER	2.4
1	B	124	THR	2.4
1	B	163	THR	2.4
1	B	170	VAL	2.4
1	B	61	LEU	2.4
1	B	251	LEU	2.4
1	B	197	PHE	2.4
1	B	144	THR	2.4
1	A	13	SER	2.3
1	A	142	VAL	2.3
1	A	26	THR	2.3
1	B	293	ILE	2.3
1	B	206	GLY	2.3
1	B	185	SER	2.3
1	B	212	SER	2.3
1	B	174	LYS	2.3
1	B	262	GLY	2.3
1	B	178	GLU	2.3
1	A	243	TYR	2.2
1	B	63	GLY	2.2
1	B	198	TYR	2.2
1	A	105	VAL	2.2
1	B	226	HIS	2.2
1	B	282	LEU	2.2
1	B	268	SER	2.2
1	B	289	ARG	2.2
1	B	177	SER	2.2
1	A	14	GLU	2.2
1	A	73	GLY	2.2
1	B	261	GLU	2.1
1	A	295	VAL	2.1
1	D	291	ASN	2.1
1	B	179	ASN	2.1
1	B	238	ILE	2.1
1	B	171	ASN	2.1
1	A	10	SER	2.0
1	B	125	SER	2.0
1	A	24	GLY	2.0
1	B	259	ARG	2.0
1	B	276	ALA	2.0
1	B	60	LYS	2.0
1	B	127	GLY	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	266	THR	2.0
1	B	243	TYR	2.0
1	A	148	PRO	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(Å ²)	Q<0.9
3	URA	B	302	8/8	0.76	0.17	38,39,44,49	0
3	URA	A	302	8/8	0.94	0.07	15,17,18,19	0
2	SO4	B	301	5/5	0.94	0.12	34,36,38,39	0
3	URA	C	302	8/8	0.95	0.06	12,15,16,17	0
3	URA	D	302	8/8	0.95	0.07	16,17,18,19	0
2	SO4	A	301	5/5	0.99	0.03	14,16,17,18	0
2	SO4	C	301	5/5	0.99	0.04	13,14,15,17	0
2	SO4	D	301	5/5	0.99	0.03	13,14,16,17	0

6.5 Other polymers [i](#)

There are no such residues in this entry.