



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2016 – 12:55 PM EDT

PDB ID : 4TXL
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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027457
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) : rb-20027457

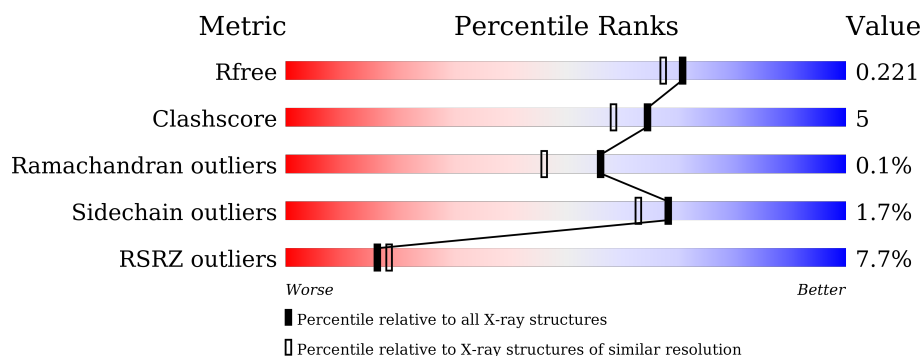
1 Overall quality at a glance

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}	91344	5832 (1.94-1.90)
Clashscore	102246	6540 (1.94-1.90)
Ramachandran outliers	100387	6464 (1.94-1.90)
Sidechain outliers	100360	6465 (1.94-1.90)
RSRZ outliers	91569	5846 (1.94-1.90)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	296	<div> <div>6%</div> <div>86% 9% 5%</div> </div>
1	B	296	<div> <div>23%</div> <div>77% 18% . .</div> </div>
1	C	296	<div> <div>92% 6% .</div> </div>
1	D	296	<div> <div>95% . . .</div> </div>

2 Entry composition [i](#)

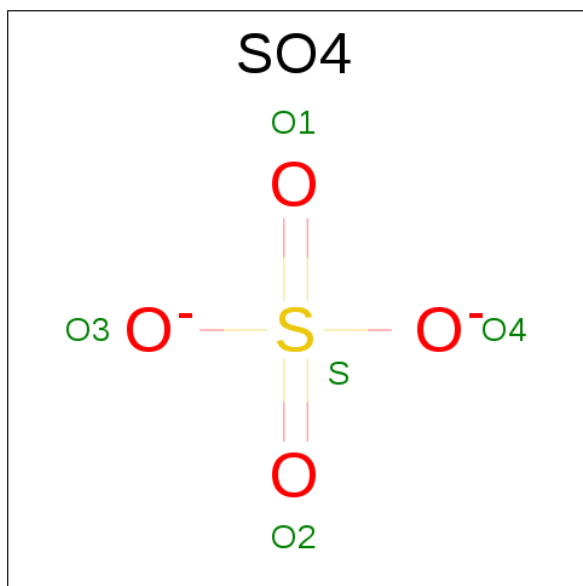
There are 4 unique types of molecules in this entry. The entry contains 9748 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 (three-letter code: SO4) (formula: O₄S).



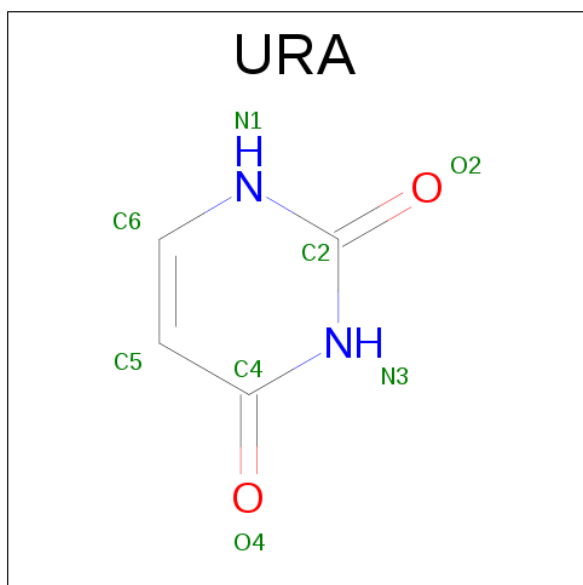
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		

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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 (three-letter code: URA) (formula: $C_4H_4N_2O_2$).



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 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

Chain C: 




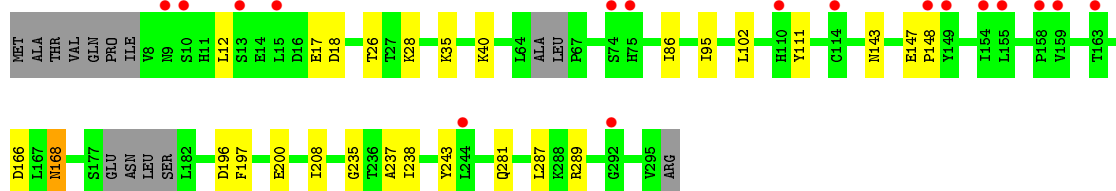
- Molecule 1: Uridine phosphorylase

Chain D: 




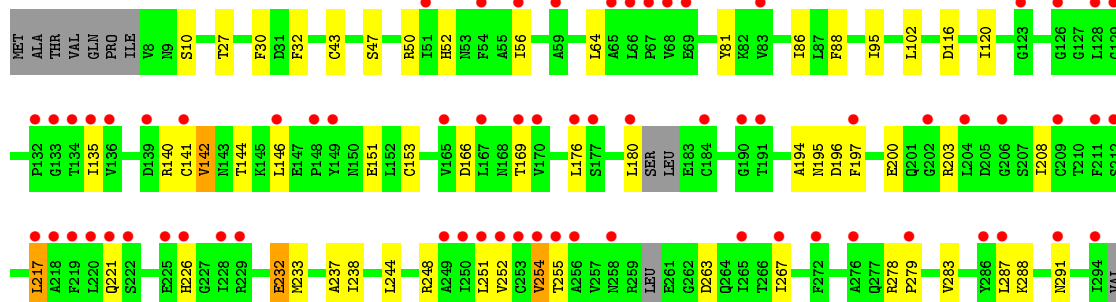
- Molecule 1: Uridine phosphorylase

Chain A: 



- Molecule 1: Uridine phosphorylase

Chain B: 



ARC

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 74.57 – 1.92	Depositor EDS
% Data completeness (in resolution range)	99.6 (71.84-1.92) 99.6 (74.57-1.92)	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 0.195 , 0.221	Depositor DCC
R_{free} test set	9006 reflections (5.03%)	DCC
Wilson B-factor (Å ²)	23.7	Xtriage
Anisotropy	0.054	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 60.9	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, 0.375 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: 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	0.25	0/2122	0.44	0/2874
1	B	0.27	0/2092	0.45	0/2839
1	C	0.29	0/2260	0.47	0/3058
1	D	0.31	0/2243	0.45	0/3035
All	All	0.28	0/8717	0.45	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	20	0
1	B	2054	0	1882	50	0
1	C	2217	0	2183	13	0
1	D	2203	0	2161	4	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	1	3
All	All	9748	0	8227	86	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 (86) 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:B:232:GLU:HB2	4:B:452:HOH:O	1.93	0.67
1:C:168:ASN:ND2	4:C:742:HOH:O	2.27	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:A:281:GLN:NE2	4:A:530:HOH:O	2.29	0.65
1:C:261:GLU:OE2	4:C:714:HOH:O	2.15	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: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:D:216:LYS:NZ	4:D:621:HOH:O	2.42	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:C:183:GLU:CD	1:C:183:GLU:H	2.14	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: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:B:47:SER:HB3	1:B:50:ARG:HB2	1.98	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: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:B:43:CYS:SG	1:B:120:ILE:HD11	2.58	0.43
1:C:95:ILE:HG23	1:C:237:ALA:HB2	2.01	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:A:196:ASP:HB3	1:A:200:GLU:HB2	2.03	0.41
1:D:47:SER:HB3	1:D:50[A]:ARG:HB2	2.02	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:A:40:LYS:HA	1:A:40:LYS:HD2	1.83	0.40
1:C:51:ILE:HG13	1:C:88:PHE:HB3	2.03	0.40
1:A:18:ASP:OD2	1:A:111:TYR:OH	2.27	0.40
1:B:86:ILE:HD13	1:B:287:LEU:HD13	2.03	0.40
1:C:216:LYS:HZ2	1:C:263:ASP:CG	2.25	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:D:439:HOH:O	4:A:405:HOH:O[1_455]	2.16	0.04
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

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%)	39	27
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%)	56	46

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%)	210 (100%)	1 (0%)	92	92
1	B	199/252 (79%)	192 (96%)	7 (4%)	43	31
1	C	236/252 (94%)	234 (99%)	2 (1%)	86	85
1	D	233/252 (92%)	227 (97%)	6 (3%)	54	44
All	All	879/1008 (87%)	863 (98%)	16 (2%)	68	60

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	30	PHE
1	C	232	GLU
1	D	30	PHE
1	D	45	CYS
1	D	50[A]	ARG
1	D	50[B]	ARG
1	D	152	LEU
1	D	232	GLU
1	A	168	ASN
1	B	30	PHE

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Mol	Chain	Res	Type
1	B	56	ILE
1	B	142	VAL
1	B	144	THR
1	B	217	LEU
1	B	232	GLU
1	B	254	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
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 carbohydrates 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	$\# Z > 2$
2	SO4	A	301	-	4,4,4	0.26	0	6,6,6	0.24	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	URA	A	302	-	6,8,8	2.97	3 (50%)	5,10,10	8.78	3 (60%)
2	SO4	B	301	-	4,4,4	0.27	0	6,6,6	0.17	0
3	URA	B	302	-	6,8,8	3.15	3 (50%)	5,10,10	8.67	3 (60%)
2	SO4	C	301	-	4,4,4	0.30	0	6,6,6	0.17	0
3	URA	C	302	-	6,8,8	2.94	3 (50%)	5,10,10	8.65	3 (60%)
2	SO4	D	301	-	4,4,4	0.24	0	6,6,6	0.11	0
3	URA	D	302	-	6,8,8	3.05	3 (50%)	5,10,10	9.33	3 (60%)

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	SO4	A	301	-	-	0/0/0/0	0/0/0/0
3	URA	A	302	-	-	0/0/0/0	0/1/1/1
2	SO4	B	301	-	-	0/0/0/0	0/0/0/0
3	URA	B	302	-	-	0/0/0/0	0/1/1/1
2	SO4	C	301	-	-	0/0/0/0	0/0/0/0
3	URA	C	302	-	-	0/0/0/0	0/1/1/1
2	SO4	D	301	-	-	0/0/0/0	0/0/0/0
3	URA	D	302	-	-	0/0/0/0	0/1/1/1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	302	URA	C2-N3	-2.85	1.32	1.38
3	D	302	URA	C2-N1	-2.79	1.32	1.38
3	B	302	URA	C2-N3	-2.73	1.32	1.38
3	D	302	URA	C2-N3	-2.71	1.32	1.38
3	A	302	URA	C2-N1	-2.66	1.32	1.38
3	B	302	URA	C2-N1	-2.61	1.32	1.38
3	A	302	URA	C2-N3	-2.55	1.32	1.38
3	C	302	URA	C2-N1	-2.51	1.33	1.38
3	C	302	URA	O4-C4	5.96	1.39	1.24
3	A	302	URA	O4-C4	6.13	1.40	1.24
3	D	302	URA	O4-C4	6.17	1.40	1.24
3	B	302	URA	O4-C4	6.52	1.41	1.24

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	302	URA	C5-C6-N1	-4.31	118.83	123.88
3	A	302	URA	C5-C6-N1	-4.07	119.12	123.88
3	B	302	URA	C5-C6-N1	-3.94	119.27	123.88
3	C	302	URA	C5-C6-N1	-3.72	119.52	123.88
3	D	302	URA	C4-N3-C2	7.20	121.80	114.21
3	A	302	URA	C4-N3-C2	7.39	121.99	114.21
3	C	302	URA	C4-N3-C2	7.45	122.06	114.21
3	B	302	URA	C4-N3-C2	7.55	122.16	114.21
3	B	302	URA	C6-N1-C2	17.30	121.23	114.47
3	C	302	URA	C6-N1-C2	17.37	121.26	114.47
3	A	302	URA	C6-N1-C2	17.64	121.36	114.47
3	D	302	URA	C6-N1-C2	19.00	121.90	114.47

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%)	0.44	17 (6%) 25 28	17, 39, 59, 72	0
1	B	284/296 (95%)	1.19	69 (24%) 1 1	17, 53, 77, 92	0
1	C	290/296 (97%)	-0.17	1 (0%) 94 94	12, 19, 33, 47	0
1	D	289/296 (97%)	-0.16	1 (0%) 94 94	12, 22, 41, 58	0
All	All	1145/1184 (96%)	0.32	88 (7%) 16 18	12, 28, 67, 92	0

All (88) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	66	LEU	6.2
1	B	149	TYR	5.4
1	B	134	THR	5.4
1	B	286	TYR	5.4
1	B	65	ALA	5.1
1	B	249	ALA	4.5
1	B	221	GLN	4.3
1	B	219	PHE	4.1
1	A	244	LEU	4.0
1	B	51	ILE	3.8
1	B	129	GLY	3.7
1	B	211	PHE	3.6
1	B	68	VAL	3.5
1	B	272	PHE	3.5
1	B	222	SER	3.5
1	B	133	GLY	3.4
1	B	67	PRO	3.4
1	B	267	ILE	3.4
1	B	83	VAL	3.3
1	B	217	LEU	3.1
1	A	155	LEU	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	180	LEU	3.1
1	B	294	ILE	3.1
1	A	75	HIS	3.1
1	B	135	ILE	3.0
1	A	13	SER	3.0
1	B	59	ALA	3.0
1	B	204	LEU	3.0
1	B	252	VAL	3.0
1	B	287	LEU	2.9
1	B	128	LEU	2.8
1	B	279	PRO	2.8
1	B	253	CYS	2.8
1	B	176	LEU	2.8
1	B	256	ALA	2.8
1	B	250	ILE	2.8
1	B	265	ILE	2.8
1	A	149	TYR	2.8
1	B	165	VAL	2.7
1	A	10	SER	2.7
1	B	54	PHE	2.7
1	A	154	ILE	2.6
1	B	56	ILE	2.6
1	B	167	LEU	2.6
1	B	148	PRO	2.6
1	B	254	VAL	2.6
1	B	258	ASN	2.6
1	B	218	ALA	2.6
1	B	191	THR	2.6
1	A	110	HIS	2.5
1	A	74	SER	2.5
1	B	220	LEU	2.5
1	B	291	ASN	2.5
1	A	15	LEU	2.5
1	A	158	PRO	2.4
1	B	255	THR	2.4
1	D	50[A]	ARG	2.4
1	C	181	SER	2.4
1	B	226	HIS	2.4
1	B	228	ILE	2.4
1	A	148	PRO	2.3
1	B	206	GLY	2.3
1	B	132	PRO	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	292	GLY	2.3
1	B	126	GLY	2.3
1	B	146	LEU	2.3
1	B	251	LEU	2.3
1	B	229	ARG	2.3
1	B	209	CYS	2.3
1	B	169	THR	2.2
1	B	139	ASP	2.2
1	B	123	GLY	2.2
1	B	202	GLY	2.2
1	A	159	VAL	2.2
1	A	114	CYS	2.2
1	B	190	GLY	2.2
1	B	276	ALA	2.2
1	B	170	VAL	2.1
1	B	184	CYS	2.1
1	B	136	VAL	2.1
1	A	9	ASN	2.1
1	B	225	GLU	2.0
1	B	212	SER	2.0
1	B	197	PHE	2.0
1	B	141	CYS	2.0
1	A	163	THR	2.0
1	B	177	SER	2.0
1	B	69	GLU	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 ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors

of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	URA	A	302	8/8	0.94	0.12	1.98	15,17,18,19	0
3	URA	D	302	8/8	0.95	0.11	0.71	16,17,18,19	0
3	URA	B	302	8/8	0.69	0.21	0.68	38,39,44,49	0
2	SO4	B	301	5/5	0.96	0.15	0.23	34,36,38,39	0
3	URA	C	302	8/8	0.96	0.10	-0.26	12,15,16,17	0
2	SO4	D	301	5/5	0.99	0.10	-0.39	13,14,16,17	0
2	SO4	C	301	5/5	0.99	0.08	-0.74	13,14,15,17	0
2	SO4	A	301	5/5	0.99	0.08	-0.87	14,16,17,18	0

6.5 Other polymers [i](#)

There are no such residues in this entry.