



Full wwPDB X-ray Structure Validation Report ⓘ

May 24, 2022 – 01:16 pm BST

PDB ID : 6Z1B
Title : Structure of K52-acetylated RutR in complex with uracil.
Authors : Kremer, M.; Schulze, S.; Lammers, M.
Deposited on : 2020-05-13
Resolution : 2.25 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.4, CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.28.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0267
CCP4 : 7.1.010 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.28.1

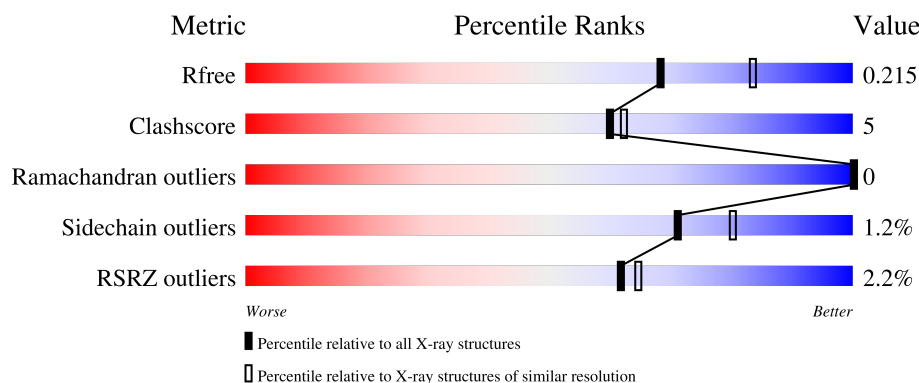
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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}	130704	1377 (2.26-2.26)
Clashscore	141614	1487 (2.26-2.26)
Ramachandran outliers	138981	1449 (2.26-2.26)
Sidechain outliers	138945	1450 (2.26-2.26)
RSRZ outliers	127900	1356 (2.26-2.26)

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	220	
1	B	220	

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
2	URA	A	301	-	X	-	-
2	URA	B	301	-	X	-	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6546 atoms, of which 3091 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.

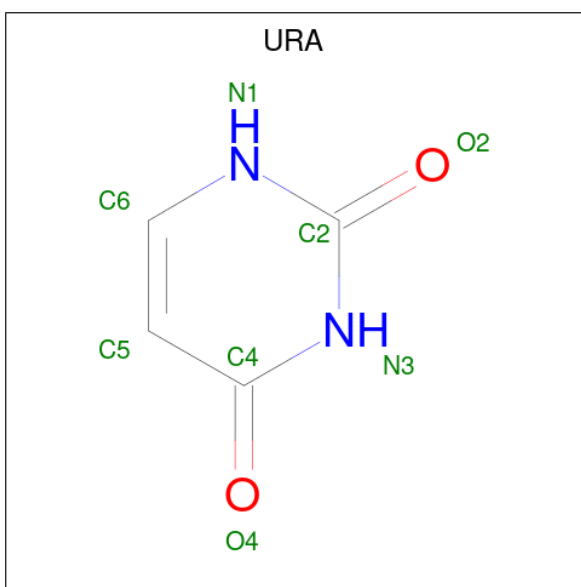
- Molecule 1 is a protein called HTH-type transcriptional regulator RutR.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	201	Total	C	H	N	O	S	0	1	0
			3134	1042	1530	272	285	5			
1	B	202	Total	C	H	N	O	S	0	1	0
			3142	1043	1535	273	286	5			

There are 16 discrepancies between the modelled and reference sequences:

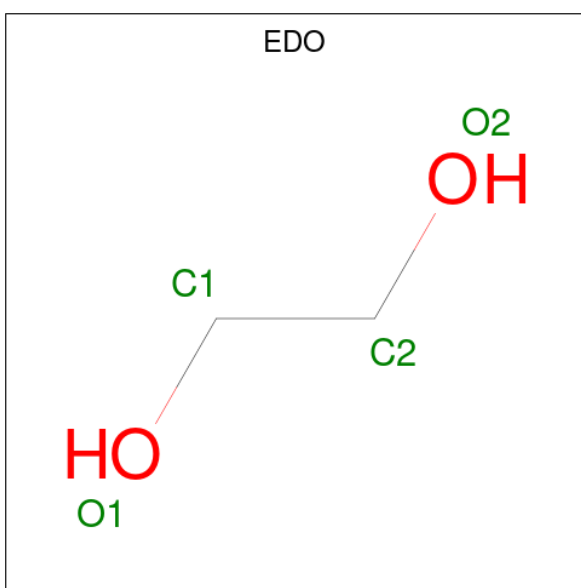
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	initiating methionine	UNP P0ACU2
A	0	GLY	-	expression tag	UNP P0ACU2
A	213	HIS	-	expression tag	UNP P0ACU2
A	214	HIS	-	expression tag	UNP P0ACU2
A	215	HIS	-	expression tag	UNP P0ACU2
A	216	HIS	-	expression tag	UNP P0ACU2
A	217	HIS	-	expression tag	UNP P0ACU2
A	218	HIS	-	expression tag	UNP P0ACU2
B	-1	MET	-	initiating methionine	UNP P0ACU2
B	0	GLY	-	expression tag	UNP P0ACU2
B	213	HIS	-	expression tag	UNP P0ACU2
B	214	HIS	-	expression tag	UNP P0ACU2
B	215	HIS	-	expression tag	UNP P0ACU2
B	216	HIS	-	expression tag	UNP P0ACU2
B	217	HIS	-	expression tag	UNP P0ACU2
B	218	HIS	-	expression tag	UNP P0ACU2

- Molecule 2 is URACIL (three-letter code: URA) (formula: C₄H₄N₂O₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	H	N	O	0	0
			12	4	4	2	2		
2	B	1	Total	C	H	N	O	0	0
			12	4	4	2	2		

- Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	H	O	0	0
			10	2	6	2		
3	A	1	Total	C	H	O	0	0
			10	2	6	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	H	O	0	0
			10	2	6	2		

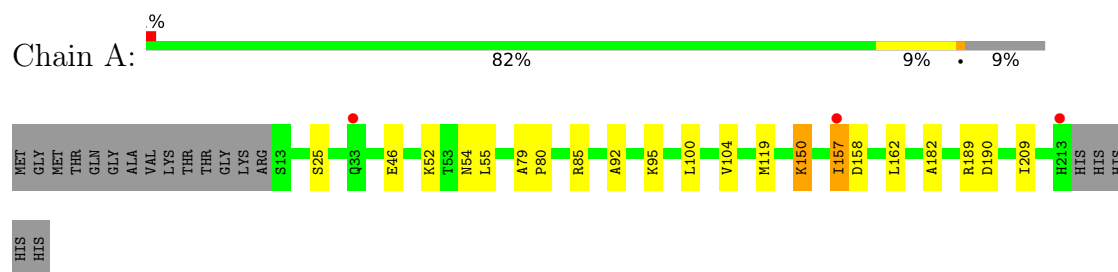
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	90	Total	O	0	0
			90	90		
4	B	126	Total	O	0	0
			126	126		

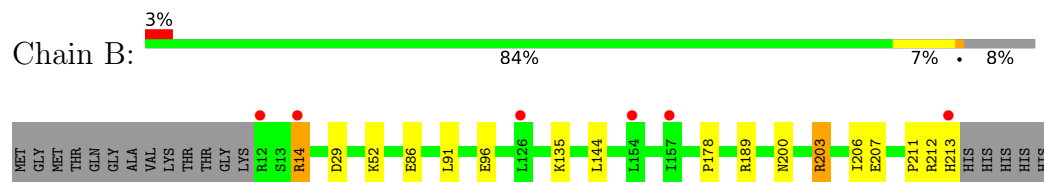
3 Residue-property plots [i](#)

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: HTH-type transcriptional regulator RutR



- Molecule 1: HTH-type transcriptional regulator RutR



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	48.31Å 88.55Å 150.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.98 – 2.25 45.98 – 2.25	Depositor EDS
% Data completeness (in resolution range)	99.9 (45.98-2.25) 99.9 (45.98-2.25)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.81 (at 2.24Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.193 , 0.215 0.193 , 0.215	Depositor DCC
R_{free} test set	1571 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	44.5	Xtriage
Anisotropy	0.269	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	(Not available) , (Not available)	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6546	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.56% 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, EDO, ALY

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	0.59	4/1629 (0.2%)	0.66	6/2205 (0.3%)
1	B	0.40	1/1637 (0.1%)	0.52	1/2216 (0.0%)
All	All	0.51	5/3266 (0.2%)	0.59	7/4421 (0.2%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	46	GLU	CB-CG	11.40	1.73	1.52
1	A	150	LYS	CE-NZ	10.20	1.74	1.49
1	A	157	ILE	CB-CG2	6.78	1.73	1.52
1	A	150	LYS	CD-CE	5.46	1.64	1.51
1	B	96	GLU	CD-OE1	5.15	1.31	1.25

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	14	ARG	CB-CG-CD	-8.27	90.10	111.60
1	A	46	GLU	OE1-CD-OE2	-7.71	114.04	123.30
1	A	150	LYS	CA-CB-CG	6.71	128.15	113.40
1	A	46	GLU	CG-CD-OE2	6.56	131.42	118.30
1	A	46	GLU	CA-CB-CG	6.55	127.82	113.40
1	A	157	ILE	C-N-CA	-5.26	108.54	121.70
1	A	150	LYS	CG-CD-CE	5.19	127.47	111.90

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	1604	1530	1644	17	1
1	B	1607	1535	1633	17	1
2	A	8	4	3	0	0
2	B	8	4	3	0	0
3	A	12	18	18	0	0
4	A	90	0	0	2	0
4	B	126	0	0	5	0
All	All	3455	3091	3301	30	1

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 (30) 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:A:150:LYS:CE	1:A:150:LYS:NZ	1.74	1.50
1:A:157:ILE:HG21	1:A:209:ILE:HG22	1.76	0.66
1:B:29:ASP:OD1	4:B:401:HOH:O	2.15	0.65
1:B:213:HIS:ND1	4:B:402:HOH:O	2.29	0.64
1:A:158:ASP:N	1:B:200:ASN:OD1	2.17	0.64
1:A:150:LYS:NZ	1:A:150:LYS:CD	2.62	0.61
1:B:212:ARG:O	1:B:213:HIS:CB	2.49	0.61
1:B:203:ARG:NH1	4:B:404:HOH:O	2.32	0.60
1:A:189:ARG:NH2	4:A:403:HOH:O	2.36	0.58
1:B:86:GLU:HA	1:B:144:LEU:HD11	1.87	0.57
1:A:157:ILE:HD12	1:A:157:ILE:O	2.06	0.55
1:A:157:ILE:HD13	1:A:162:LEU:HB2	1.92	0.51
1:B:86:GLU:CA	1:B:144:LEU:HD11	2.40	0.51
1:A:190:ASP:OD2	4:A:401:HOH:O	2.20	0.49
1:A:119:MET:HE1	1:B:178:PRO:HB2	1.95	0.48
1:B:203:ARG:HG3	4:B:403:HOH:O	2.13	0.47
1:A:157:ILE:HG22	1:B:203:ARG:HD2	1.97	0.46
1:B:211:PRO:O	1:B:212:ARG:HB2	2.16	0.46
1:B:91:LEU:HD22	1:B:206:ILE:HG23	1.98	0.45
1:A:190:ASP:OD1	1:A:190:ASP:C	2.55	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:92:ALA:HA	1:A:95:LYS:HG2	2.00	0.43
1:A:79:ALA:HB3	1:A:80:PRO:HD3	2.01	0.42
1:B:212:ARG:HD3	1:B:212:ARG:HA	1.70	0.42
1:A:119:MET:HE3	1:A:119:MET:HA	2.01	0.42
1:B:203:ARG:NH2	1:B:207:GLU:HG3	2.34	0.42
1:A:55:LEU:HD23	1:A:55:LEU:C	2.40	0.42
1:B:189:ARG:NH1	4:B:414:HOH:O	2.53	0.41
1:A:100:LEU:O	1:A:104:VAL:HG23	2.21	0.41
1:A:182:ALA:O	1:B:135:LYS:HD2	2.20	0.41
1:B:212:ARG:O	1:B:213:HIS:HB3	2.20	0.40

All (1) 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)
1:A:157:ILE:O	1:B:14:ARG:NH2[3_454]	1.89	0.31

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	199/220 (90%)	198 (100%)	1 (0%)	0	100	100
1	B	200/220 (91%)	198 (99%)	2 (1%)	0	100	100
All	All	399/440 (91%)	396 (99%)	3 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	166/180 (92%)	163 (98%)	3 (2%)	59	68
1	B	167/180 (93%)	166 (99%)	1 (1%)	86	91
All	All	333/360 (92%)	329 (99%)	4 (1%)	71	80

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	SER
1	A	54	ASN
1	A	85	ARG
1	B	203	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	33	GLN
1	A	43	GLN
1	A	179	GLN
1	A	195	ASN
1	A	196	GLN
1	A	213	HIS
1	B	33	GLN
1	B	179	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	ALY	B	52	1	10,11,12	0.85	0	7,12,14	1.44	1 (14%)
1	ALY	A	52	1	10,11,12	0.81	0	7,12,14	1.68	3 (42%)

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
1	ALY	B	52	1	-	4/9/10/12	-
1	ALY	A	52	1	-	3/9/10/12	-

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	52	ALY	CE-NZ-CH	2.64	126.61	122.56
1	B	52	ALY	CE-NZ-CH	2.43	126.29	122.56
1	A	52	ALY	CD-CE-NZ	-2.35	105.49	112.21
1	A	52	ALY	CD-CG-CB	-2.18	105.92	113.62

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	52	ALY	N-CA-CB-CG
1	A	52	ALY	C-CA-CB-CG
1	B	52	ALY	N-CA-CB-CG
1	B	52	ALY	C-CA-CB-CG
1	B	52	ALY	CD-CE-NZ-CH
1	A	52	ALY	CG-CD-CE-NZ
1	B	52	ALY	CG-CD-CE-NZ

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

5 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	URA	A	301	-	6,8,8	4.83	6 (100%)	4,10,10	7.51	3 (75%)
3	EDO	A	302	-	3,3,3	0.48	0	2,2,2	0.24	0
3	EDO	A	304	-	3,3,3	0.50	0	2,2,2	0.30	0
2	URA	B	301	-	6,8,8	4.90	6 (100%)	4,10,10	7.46	3 (75%)
3	EDO	A	303	-	3,3,3	0.49	0	2,2,2	0.20	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
2	URA	A	301	-	-	-	0/1/1/1
3	EDO	A	302	-	-	0/1/1/1	-
3	EDO	A	304	-	-	0/1/1/1	-
2	URA	B	301	-	-	-	0/1/1/1
3	EDO	A	303	-	-	1/1/1/1	-

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	301	URA	C4-N3	7.14	1.45	1.33
2	B	301	URA	C4-N3	7.05	1.45	1.33
2	B	301	URA	C2-N1	5.42	1.48	1.38
2	A	301	URA	C2-N1	5.26	1.48	1.38
2	B	301	URA	C2-N3	5.20	1.48	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	301	URA	C6-N1	5.18	1.45	1.34
2	A	301	URA	C2-N3	5.05	1.48	1.38
2	A	301	URA	C6-N1	4.93	1.45	1.34
2	A	301	URA	C5-C6	2.65	1.43	1.38
2	B	301	URA	C5-C6	2.58	1.43	1.38
2	A	301	URA	O4-C4	-2.16	1.19	1.24
2	B	301	URA	O4-C4	-2.09	1.19	1.24

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	URA	N3-C2-N1	-10.67	119.95	128.43
2	B	301	URA	N3-C2-N1	-10.63	119.98	128.43
2	A	301	URA	C6-N1-C2	10.20	119.45	114.42
2	B	301	URA	C6-N1-C2	10.08	119.39	114.42
2	B	301	URA	C5-C6-N1	-2.38	121.00	123.96
2	A	301	URA	C5-C6-N1	-2.25	121.17	123.96

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	303	EDO	O1-C1-C2-O2

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	200/220 (90%)	0.17	3 (1%) 73 75	35, 53, 84, 121	0
1	B	201/220 (91%)	0.14	6 (2%) 50 53	34, 53, 78, 123	0
All	All	401/440 (91%)	0.15	9 (2%) 62 65	34, 53, 81, 123	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	213	HIS	4.5
1	A	157	ILE	2.9
1	B	14	ARG	2.9
1	B	12	ARG	2.8
1	B	154	LEU	2.3
1	A	213	HIS	2.3
1	A	33	GLN	2.2
1	B	157	ILE	2.1
1	B	126	LEU	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	ALY	B	52	12/13	0.86	0.16	75,99,133,133	0
1	ALY	A	52	12/13	0.92	0.19	49,75,108,108	0

6.3 Carbohydrates [i](#)

There are no monosaccharides 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(\AA^2)	Q<0.9
3	EDO	A	304	4/4	0.63	0.15	90,108,111,113	0
3	EDO	A	303	4/4	0.83	0.13	75,90,91,92	0
3	EDO	A	302	4/4	0.93	0.30	90,108,112,113	0
2	URA	B	301	8/8	0.94	0.16	47,49,59,59	0
2	URA	A	301	8/8	0.95	0.17	41,45,52,54	0

6.5 Other polymers [i](#)

There are no such residues in this entry.