



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 06:34 am GMT

PDB ID : 3QE7  
Title : Crystal Structure of Uracil Transporter–UraA  
Authors : Lu, F.R.; Li, S.; Yan, N.  
Deposited on : 2011-01-20  
Resolution : 2.78 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](mailto: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.

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

MolProbity : 4.02b-467  
Mogul : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
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) : recalc28949

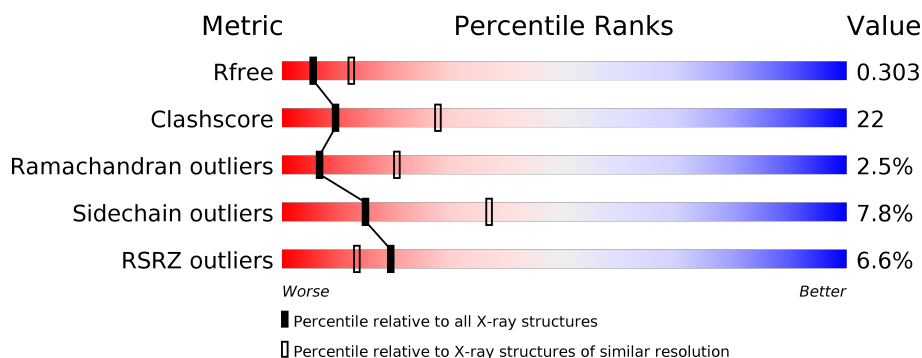
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.78 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))
$R_{free}$	100719	3276 (2.80-2.76)
Clashscore	112137	3771 (2.80-2.76)
Ramachandran outliers	110173	3707 (2.80-2.76)
Sidechain outliers	110143	3709 (2.80-2.76)
RSRZ outliers	101464	3307 (2.80-2.76)

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  $\geq 3$ , 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	429	<div> <div>6%</div> <div>57%</div> <div>32%</div> <div>5%</div> <div>5%</div> </div>

## 2 Entry composition [i](#)

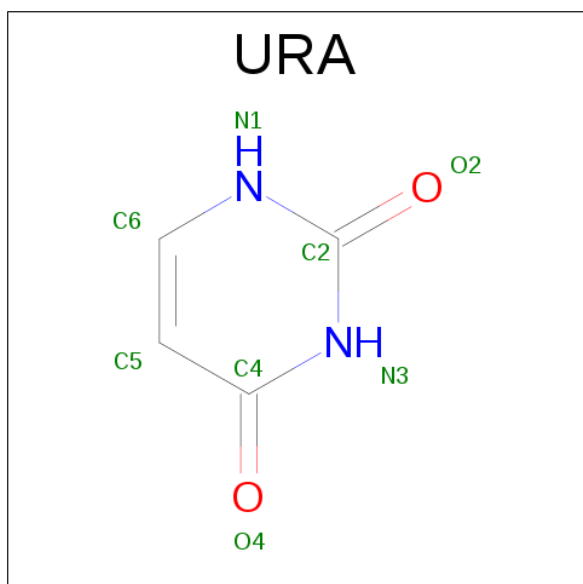
There are 3 unique types of molecules in this entry. The entry contains 3030 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Uracil permease.

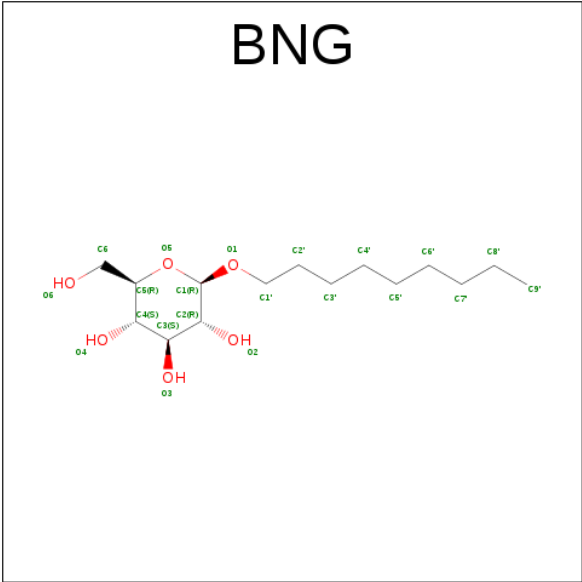
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	407	3001	2004	475	508	14	0	0	0

- Molecule 2 is URACIL (three-letter code: URA) (formula: C<sub>4</sub>H<sub>4</sub>N<sub>2</sub>O<sub>2</sub>).



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

- Molecule 3 is SUGAR (B-NONYLGLUCOSIDE) (three-letter code: BNG) (formula: C<sub>15</sub>H<sub>30</sub>O<sub>6</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			21	15	6		



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 64 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	96.76Å 96.76Å 251.95Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	41.92 – 2.78 41.92 – 2.78	Depositor EDS
% Data completeness (in resolution range)	99.0 (41.92-2.78) 99.1 (41.92-2.78)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.07 (at 2.77Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.3_473)	Depositor
R, $R_{free}$	0.248 , 0.303 0.244 , 0.303	Depositor DCC
$R_{free}$ test set	929 reflections (5.11%)	DCC
Wilson B-factor (Å <sup>2</sup> )	71.3	Xtriage
Anisotropy	0.403	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 78.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	3030	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	77.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.01% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: BNG, URA

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	0.28	0/3066	0.49	1/4187 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	39	PHE	N-CA-C	5.55	125.99	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	260	ASP	Peptide

### 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	3001	0	3224	137	0
2	A	8	0	3	0	0
3	A	21	0	29	2	0
All	All	3030	0	3256	137	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 22.

All (137) 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:179:PHE:O	1:A:183:ILE:HB	1.68	0.94
1:A:50:ASN:O	1:A:54:THR:HG23	1.71	0.90
1:A:33:VAL:HG23	1:A:43:PRO:HB2	1.57	0.85
1:A:12:PRO:HB2	1:A:17:THR:HG22	1.58	0.82
1:A:260:ASP:HB3	1:A:261:PRO:HD2	1.65	0.78
1:A:64:LYS:HD3	1:A:261:PRO:HD3	1.65	0.78
1:A:386:ALA:HB1	1:A:387:ALA:HA	1.64	0.77
1:A:238:VAL:HG13	1:A:286:THR:HG23	1.65	0.77
1:A:260:ASP:O	1:A:262:GLY:N	2.19	0.76
1:A:54:THR:HG22	1:A:69:LEU:H	1.52	0.75
1:A:64:LYS:HB3	1:A:261:PRO:HG3	1.69	0.74
1:A:274:THR:CG2	1:A:285:ASN:HB2	2.17	0.74
1:A:360:TYR:OH	1:A:369:THR:HG21	1.90	0.72
1:A:214:ALA:O	1:A:318:CYS:HB2	1.92	0.68
1:A:42:ASN:O	1:A:45:THR:HG23	1.92	0.68
1:A:350:ILE:HD11	3:A:431:BNG:H5'1	1.75	0.67
1:A:42:ASN:HD22	1:A:217:THR:CG2	2.08	0.67
1:A:224:GLU:O	1:A:228:ILE:HG12	1.95	0.67
1:A:33:VAL:HG21	1:A:326:ILE:HG21	1.75	0.67
1:A:42:ASN:HD22	1:A:217:THR:HG22	1.58	0.66
1:A:54:THR:CG2	1:A:69:LEU:H	2.08	0.66
1:A:156:SER:O	1:A:160:ILE:HG12	1.97	0.65
1:A:226:PHE:O	1:A:230:THR:HG23	1.98	0.64
1:A:6:ILE:HD12	1:A:300:VAL:HG22	1.79	0.64
1:A:283:THR:HG23	1:A:284:PRO:HD2	1.80	0.62
1:A:64:LYS:CB	1:A:261:PRO:HG3	2.29	0.61
1:A:274:THR:HG22	1:A:285:ASN:HB2	1.83	0.61
1:A:216:PRO:HD3	1:A:318:CYS:HB3	1.83	0.60
1:A:34:LEU:HD23	1:A:334:MET:HE3	1.82	0.60
1:A:157:LYS:O	1:A:161:ILE:HG13	2.02	0.59
1:A:127:ALA:HB2	1:A:369:THR:HB	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:379:GLY:CA	1:A:380:ALA:HB3	2.34	0.58
1:A:205:THR:HB	1:A:206:PRO:HD3	1.85	0.57
1:A:386:ALA:CB	1:A:387:ALA:HA	2.29	0.57
1:A:28:MET:HE3	1:A:68:TYR:HE2	1.68	0.56
1:A:78:PRO:HB2	1:A:231:ILE:HD13	1.86	0.56
1:A:260:ASP:HB3	1:A:261:PRO:CD	2.35	0.56
1:A:54:THR:HG22	1:A:69:LEU:HB2	1.87	0.56
1:A:34:LEU:HA	1:A:334:MET:CE	2.36	0.56
1:A:183:ILE:O	1:A:187:ILE:HG13	2.05	0.55
1:A:351:ARG:O	1:A:355:GLU:HG2	2.06	0.55
1:A:162:SER:O	1:A:166:LEU:HG	2.07	0.54
1:A:183:ILE:N	1:A:184:PRO:HD2	2.21	0.54
1:A:18:ILE:N	1:A:19:PRO:HD2	2.22	0.54
1:A:123:ALA:O	1:A:369:THR:HG22	2.07	0.54
1:A:34:LEU:HA	1:A:334:MET:HE3	1.90	0.54
1:A:358:VAL:HB	1:A:360:TYR:CE2	2.43	0.54
1:A:379:GLY:HA2	1:A:380:ALA:HB3	1.89	0.53
1:A:33:VAL:CG2	1:A:326:ILE:HG21	2.38	0.53
1:A:4:ARG:O	1:A:5:ALA:HB2	2.10	0.52
1:A:207:ILE:O	1:A:208:ILE:HG13	2.10	0.52
1:A:347:ALA:HA	1:A:350:ILE:HD11	1.91	0.52
1:A:180:LEU:O	1:A:180:LEU:HD13	2.10	0.52
1:A:162:SER:O	1:A:165:THR:HG22	2.10	0.51
1:A:183:ILE:HG22	1:A:184:PRO:HD3	1.92	0.51
1:A:183:ILE:N	1:A:184:PRO:CD	2.72	0.51
1:A:401:GLY:O	1:A:405:ILE:HG12	2.09	0.51
1:A:226:PHE:CD1	1:A:227:ALA:N	2.79	0.51
1:A:25:LEU:HD21	1:A:316:LEU:HD12	1.93	0.51
1:A:67:ALA:O	1:A:306:ILE:HG21	2.10	0.51
1:A:373:LEU:O	1:A:377:VAL:HG12	2.11	0.51
1:A:33:VAL:HG23	1:A:43:PRO:CB	2.36	0.50
1:A:120:PHE:CZ	1:A:400:ILE:HD11	2.46	0.50
1:A:201:ILE:HG22	1:A:207:ILE:HG13	1.92	0.50
1:A:360:TYR:CE1	1:A:369:THR:HG21	2.47	0.50
1:A:365:ASN:O	1:A:369:THR:HG23	2.12	0.50
1:A:39:PHE:O	1:A:83:LEU:HD11	2.12	0.50
1:A:161:ILE:O	1:A:165:THR:HG22	2.11	0.50
1:A:316:LEU:O	1:A:319:VAL:HB	2.12	0.49
1:A:179:PHE:CG	1:A:180:LEU:N	2.80	0.49
1:A:33:VAL:HG21	1:A:326:ILE:CG2	2.43	0.49
1:A:371:VAL:O	1:A:375:ILE:HG12	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:160:ILE:O	1:A:164:THR:HG23	2.12	0.49
1:A:263:LEU:HG	1:A:267:MET:HE2	1.95	0.49
1:A:180:LEU:HA	1:A:183:ILE:HG22	1.94	0.49
1:A:224:GLU:HB3	1:A:226:PHE:CE2	2.48	0.49
1:A:141:ALA:HB1	1:A:147:LEU:HG	1.95	0.48
1:A:360:TYR:HE1	1:A:369:THR:CG2	2.25	0.48
1:A:251:ASN:HD21	1:A:357:LYS:HE3	1.79	0.48
1:A:274:THR:HG21	1:A:285:ASN:H	1.78	0.48
1:A:186:LEU:O	1:A:190:LEU:HG	2.13	0.48
1:A:199:MET:O	1:A:202:VAL:HG12	2.13	0.48
1:A:385:GLY:HA2	1:A:386:ALA:C	2.34	0.48
1:A:8:VAL:HG13	1:A:303:THR:HG23	1.96	0.47
1:A:119:LEU:HD22	1:A:400:ILE:HG23	1.96	0.47
1:A:113:THR:HB	1:A:116:LEU:HD12	1.97	0.47
1:A:226:PHE:CG	1:A:227:ALA:N	2.83	0.47
1:A:379:GLY:HA2	1:A:380:ALA:O	2.15	0.47
1:A:160:ILE:O	1:A:164:THR:CG2	2.63	0.47
1:A:101:PHE:HB3	1:A:271:GLY:HA3	1.97	0.46
1:A:150:GLU:HA	1:A:151:GLY:HA2	1.50	0.46
1:A:207:ILE:O	1:A:207:ILE:HG22	2.16	0.46
1:A:393:ALA:O	1:A:397:ILE:HG12	2.16	0.46
1:A:65:ILE:HG12	1:A:257:LEU:HD22	1.97	0.46
1:A:64:LYS:CD	1:A:261:PRO:HD3	2.43	0.46
1:A:23:GLN:HE22	1:A:298:THR:HG21	1.80	0.46
1:A:381:LYS:HG2	1:A:390:LYS:HG2	1.98	0.46
1:A:171:LEU:HA	1:A:174:VAL:HG12	1.98	0.45
1:A:87:TYR:CZ	1:A:91:LEU:HD21	2.51	0.45
1:A:212:TRP:O	1:A:319:VAL:HA	2.17	0.44
1:A:24:HIS:CD2	1:A:291:ASN:HD22	2.35	0.44
1:A:318:CYS:O	1:A:318:CYS:SG	2.76	0.44
1:A:54:THR:HG22	1:A:69:LEU:CB	2.48	0.44
1:A:11:ARG:HB3	1:A:304:TRP:CD2	2.53	0.44
1:A:66:PRO:HB2	1:A:295:MET:SD	2.57	0.44
1:A:94:PHE:CG	1:A:284:PRO:HD3	2.53	0.44
1:A:18:ILE:HD12	1:A:312:PHE:HZ	1.83	0.43
1:A:290:GLU:HA	1:A:345:ILE:HD13	2.00	0.43
1:A:379:GLY:N	1:A:380:ALA:HB3	2.32	0.43
1:A:89:VAL:HG12	1:A:224:GLU:HG3	1.99	0.43
1:A:57:TYR:CE2	1:A:67:ALA:HB3	2.54	0.43
1:A:350:ILE:O	1:A:354:ILE:HG13	2.18	0.43
1:A:360:TYR:CZ	1:A:369:THR:HG21	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:39:PHE:N	1:A:40:HIS:HA	2.33	0.43
1:A:207:ILE:HD11	1:A:325:ALA:N	2.33	0.43
1:A:144:ALA:HA	1:A:331:LEU:HG	2.01	0.42
1:A:14:LEU:O	1:A:18:ILE:HG12	2.18	0.42
1:A:179:PHE:O	1:A:183:ILE:CB	2.54	0.42
1:A:360:TYR:HE1	1:A:369:THR:HG23	1.84	0.42
1:A:388:GLU:HG3	1:A:390:LYS:HG3	2.01	0.42
1:A:83:LEU:HA	1:A:83:LEU:HD23	1.84	0.42
1:A:14:LEU:HD11	1:A:18:ILE:HD11	2.02	0.42
1:A:56:LEU:HD21	1:A:272:LEU:HD12	2.01	0.42
1:A:120:PHE:HZ	1:A:400:ILE:HD11	1.83	0.42
1:A:201:ILE:HG22	1:A:207:ILE:CG1	2.50	0.41
1:A:34:LEU:HD23	1:A:334:MET:CE	2.49	0.41
1:A:162:SER:HA	1:A:165:THR:HG22	2.02	0.41
1:A:289:GLY:HA3	3:A:431:BNG:O4	2.20	0.41
1:A:172:GLY:HA2	1:A:176:PHE:CE2	2.56	0.41
1:A:77:SER:N	1:A:78:PRO:HD2	2.35	0.41
1:A:13:PRO:O	1:A:17:THR:HG23	2.21	0.41
1:A:179:PHE:CE1	1:A:180:LEU:HB2	2.56	0.41
1:A:407:LYS:HA	1:A:407:LYS:HD2	1.95	0.41
1:A:56:LEU:HD21	1:A:272:LEU:CD1	2.51	0.40
1:A:337:VAL:O	1:A:340:LEU:HB3	2.22	0.40
1:A:65:ILE:HD11	1:A:261:PRO:HB2	2.04	0.40
1:A:363:ALA:O	1:A:367:ILE:HG12	2.22	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	405/429 (94%)	364 (90%)	31 (8%)	10 (2%)	6	20

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	5	ALA
1	A	261	PRO
1	A	391	GLY
1	A	208	ILE
1	A	379	GLY
1	A	318	CYS
1	A	389	LEU
1	A	260	ASP
1	A	375	ILE
1	A	319	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	320/340 (94%)	295 (92%)	25 (8%)	15	37

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	ARG
1	A	34	LEU
1	A	45	THR
1	A	82	LEU
1	A	89	VAL
1	A	101	PHE
1	A	147	LEU
1	A	153	THR
1	A	164	THR
1	A	176	PHE
1	A	217	THR
1	A	220	THR
1	A	232	LEU
1	A	246	LEU
1	A	270	ASN

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Mol	Chain	Res	Type
1	A	272	LEU
1	A	286	THR
1	A	288	TYR
1	A	304	TRP
1	A	319	VAL
1	A	339	LEU
1	A	340	LEU
1	A	350	ILE
1	A	362	LYS
1	A	369	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	16	GLN
1	A	23	GLN
1	A	24	HIS
1	A	40	HIS
1	A	251	ASN

### 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](#)

2 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	430	-	6,8,8	1.15	0	5,10,10	8.34	4 (80%)
3	BNG	A	431	-	21,21,21	1.13	2 (9%)	26,26,26	1.44	3 (11%)

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	430	-	-	0/0/0/0	0/1/1/1
3	BNG	A	431	-	-	0/12/32/32	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	431	BNG	O4-C4	-2.84	1.36	1.43
3	A	431	BNG	O2-C2	-2.41	1.37	1.43

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	430	URA	N1-C2-N3	-10.22	121.05	128.40
2	A	430	URA	C5-C6-N1	-3.88	119.45	123.92
3	A	431	BNG	C1-O5-C5	-2.79	108.46	113.72
3	A	431	BNG	C4-C3-C2	-2.72	106.04	110.84
3	A	431	BNG	C1'-O1-C1	4.90	122.27	113.87
2	A	430	URA	C4-N3-C2	7.98	120.99	114.13
2	A	430	URA	C6-N1-C2	12.73	120.56	114.41

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	431	BNG	2	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	407/429 (94%)	0.06	27 (6%) 19 13	42, 66, 144, 214	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	383	ASN	6.4
1	A	384	ILE	5.9
1	A	385	GLY	5.9
1	A	199	MET	4.7
1	A	185	ILE	4.1
1	A	179	PHE	3.9
1	A	150	GLU	3.7
1	A	176	PHE	3.5
1	A	207	ILE	3.2
1	A	202	VAL	3.2
1	A	181	ALA	3.0
1	A	110	LYS	2.9
1	A	211	HIS	2.8
1	A	62	LYS	2.7
1	A	209	ASN	2.7
1	A	177	ARG	2.7
1	A	193	TYR	2.7
1	A	151	GLY	2.6
1	A	115	TRP	2.6
1	A	255	LYS	2.4
1	A	260	ASP	2.3
1	A	156	SER	2.2
1	A	379	GLY	2.1
1	A	381	LYS	2.1
1	A	180	LEU	2.1
1	A	386	ALA	2.1
1	A	213	PHE	2.1



## 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 carbohydrates 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. 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, 95<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
3	BNG	A	431	21/21	0.94	0.24	1.90	43,56,81,82	0
2	URA	A	430	8/8	0.96	0.14	-1.26	39,41,51,58	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.