



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:58 PM GMT

PDB ID : 1I4S  
Title : CRYSTAL STRUCTURE OF RNASE III ENDONUCLEASE DOMAIN  
FROM AQUIFEX AEOLICUS AT 2.15 ANGSTROM RESOLUTION  
Authors : Blaszczyk, J.; Ji, X.  
Deposited on : 2001-02-22  
Resolution : 2.15 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 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

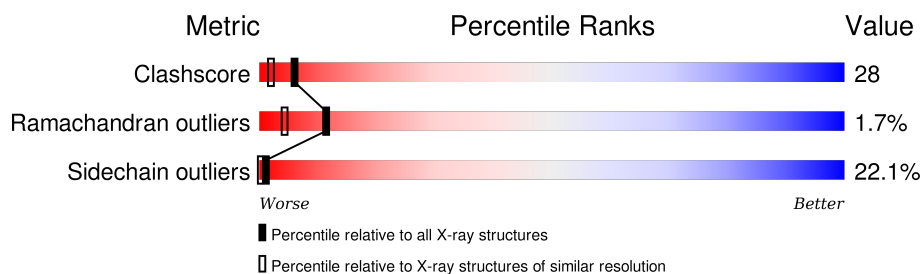
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.15 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))
Clashscore	102246	1152 (2.16-2.16)
Ramachandran outliers	100387	1131 (2.16-2.16)
Sidechain outliers	100360	1131 (2.16-2.16)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	147	
1	B	147	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2696 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 RIBONUCLEASE III.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	147	Total	C	N	O	S	0	0	0
			1222	803	197	220	2			
1	B	147	Total	C	N	O	S	0	0	0
			1222	803	197	220	2			

- Molecule 2 is water.

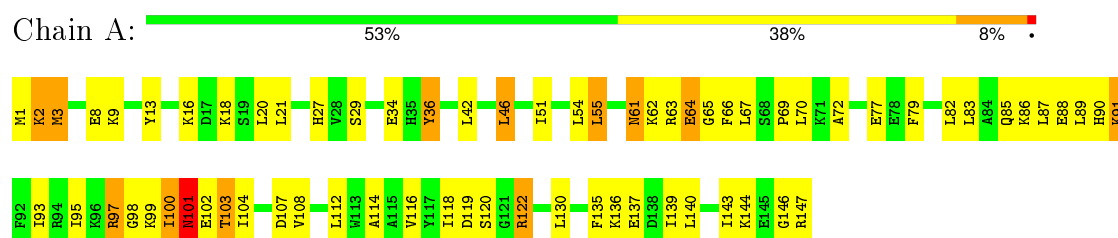
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	173	Total	O	0	0
			173	173		
2	B	79	Total	O	0	0
			79	79		

### 3 Residue-property plots

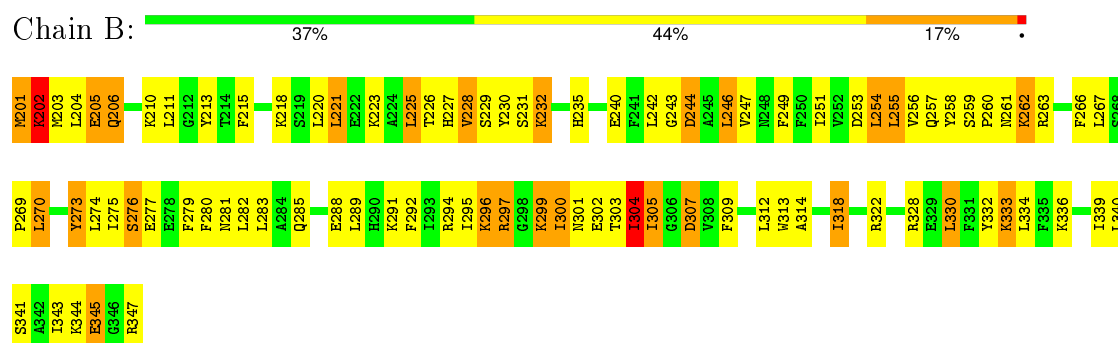
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

Note EDS was not executed.

#### • Molecule 1: RIBONUCLEASE III



#### • Molecule 1: RIBONUCLEASE III



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	46.16 Å 51.19 Å 123.47 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.15	Depositor
% Data completeness (in resolution range)	94.8 (30.00-2.15)	Depositor
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	SHELXL-97	Depositor
R, $R_{free}$	0.212 , 0.275	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2696	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	54.0	wwPDB-VP

## 5 Model quality

### 5.1 Standard geometry

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	0/1247	0.90	2/1670 (0.1%)
1	B	0.32	0/1247	0.83	1/1670 (0.1%)
All	All	0.32	0/2494	0.87	3/3340 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	244	ASP	CB-CG-OD1	6.79	124.41	118.30
1	A	122	ARG	NE-CZ-NH1	-6.41	117.10	120.30
1	A	36	TYR	CB-CG-CD1	6.13	124.68	121.00

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	1222	0	1254	49	0
1	B	1222	0	1251	94	0
2	A	173	0	0	2	0
2	B	79	0	0	2	0
All	All	2696	0	2505	141	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 28.

All (141) 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:228:VAL:HG12	1:B:232:LYS:HA	1.43	0.97
1:B:336:LYS:HE2	1:B:340:LEU:HD11	1.51	0.92
1:A:93:ILE:HG12	1:A:108:VAL:HG21	1.54	0.89
1:B:229:SER:HB2	1:B:296:LYS:HB3	1.57	0.86
1:B:295:ILE:HD11	1:B:299:LYS:HD3	1.65	0.78
1:A:82:LEU:O	1:A:86:LYS:HE2	1.85	0.77
1:A:20:LEU:HD12	1:A:119:ASP:HB2	1.67	0.76
1:B:229:SER:OG	1:B:295:ILE:HD12	1.85	0.76
1:B:266:PHE:O	1:B:269:PRO:HD2	1.85	0.76
1:B:277:GLU:OE1	1:B:302:GLU:HG2	1.86	0.74
1:B:261:ASN:HD22	1:B:266:PHE:HB2	1.54	0.73
1:A:101:ASN:HD22	1:A:101:ASN:N	1.87	0.73
1:B:210:LYS:HE3	1:B:288:GLU:HG3	1.73	0.69
1:A:90:HIS:NE2	1:A:91:LYS:HE3	2.09	0.68
1:B:288:GLU:O	1:B:291:LYS:HG2	1.92	0.68
1:A:100:ILE:O	1:A:104:ILE:HB	1.94	0.68
1:B:201:MET:HB3	1:B:202:LYS:HD3	1.76	0.66
1:A:90:HIS:CE1	1:A:91:LYS:HE3	2.30	0.66
1:B:330:LEU:O	1:B:334:LEU:HG	1.96	0.65
1:A:82:LEU:HD22	1:A:85:GLN:NE2	2.11	0.65
1:B:260:PRO:HD2	1:B:266:PHE:CE2	2.33	0.64
1:A:139:ILE:O	1:A:143:ILE:HG13	1.97	0.64
1:A:8:GLU:HB3	1:A:13:TYR:O	1.99	0.63
1:B:202:LYS:N	1:B:202:LYS:HD3	2.14	0.62
1:B:255:LEU:HD21	1:B:270:LEU:HD13	1.81	0.62
1:B:261:ASN:ND2	1:B:263:ARG:HB2	2.15	0.62
1:A:1:MET:HG2	1:A:2:LYS:H	1.66	0.60
1:B:259:SER:O	1:B:262:LYS:HE3	2.01	0.60
1:B:201:MET:C	1:B:202:LYS:HD3	2.22	0.60
1:A:34:GLU:HA	1:A:34:GLU:OE1	2.02	0.60
1:B:261:ASN:HD22	1:B:266:PHE:CB	2.15	0.60
1:B:254:LEU:HD13	1:B:332:TYR:CZ	2.38	0.59
1:B:300:ILE:O	1:B:304:ILE:HD12	2.03	0.59
1:B:213:TYR:HB2	1:B:330:LEU:HD11	1.85	0.58
1:B:228:VAL:HG13	1:B:235:HIS:CD2	2.38	0.58
1:B:261:ASN:HB3	1:B:266:PHE:CD2	2.39	0.57
1:B:260:PRO:HD2	1:B:266:PHE:HE2	1.68	0.57
1:B:281:ASN:OD1	1:B:305:ILE:HG12	2.04	0.57
1:A:46:LEU:HD23	1:B:249:PHE:CE1	2.40	0.57
1:B:215:PHE:CD1	1:B:221:LEU:HG	2.40	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:89:LEU:HD21	1:A:112:LEU:HD22	1.86	0.56
1:B:301:ASN:HD22	1:B:304:ILE:HD12	1.71	0.56
1:B:270:LEU:HD11	1:B:343:ILE:HG23	1.86	0.56
1:B:261:ASN:HB3	1:B:266:PHE:CG	2.41	0.56
1:B:225:LEU:HD23	1:B:289:LEU:HD21	1.86	0.56
1:A:87:LEU:O	1:A:88:GLU:HG2	2.05	0.55
1:B:229:SER:O	1:B:296:LYS:HE2	2.06	0.54
1:B:295:ILE:HD11	1:B:299:LYS:CD	2.37	0.54
1:B:273:TYR:O	1:B:276:SER:HB3	2.08	0.53
1:B:336:LYS:HE2	1:B:340:LEU:CD1	2.32	0.53
1:A:77:GLU:OE1	1:A:103:THR:HB	2.08	0.53
1:B:263:ARG:O	1:B:267:LEU:HG	2.08	0.53
1:B:254:LEU:HD23	1:B:339:ILE:HD13	1.91	0.53
1:B:210:LYS:HE3	1:B:288:GLU:CG	2.39	0.53
1:A:136:LYS:HE3	1:A:140:LEU:HD11	1.90	0.52
1:B:303:THR:HG22	1:B:303:THR:O	2.10	0.52
1:A:63:ARG:O	1:A:67:LEU:HG	2.09	0.52
1:B:301:ASN:O	1:B:304:ILE:HB	2.10	0.52
1:B:255:LEU:CD2	1:B:270:LEU:HD13	2.40	0.52
1:A:101:ASN:N	1:A:101:ASN:ND2	2.55	0.52
1:A:79:PHE:CE2	1:A:83:LEU:HD11	2.46	0.51
1:A:114:ALA:O	1:A:118:ILE:HG13	2.10	0.51
1:A:82:LEU:HD22	1:A:85:GLN:HE22	1.76	0.51
1:B:277:GLU:OE2	1:B:302:GLU:O	2.29	0.51
1:B:299:LYS:O	1:B:301:ASN:ND2	2.45	0.50
1:B:279:PHE:CE2	1:B:283:LEU:HD21	2.47	0.50
1:B:228:VAL:HG13	1:B:235:HIS:CE1	2.47	0.50
1:B:275:ILE:HG22	1:B:275:ILE:O	2.12	0.50
1:A:29:SER:OG	1:A:95:ILE:HG13	2.11	0.50
1:B:328:ARG:HD3	1:B:332:TYR:CE2	2.46	0.49
1:A:64:GLU:HG3	1:A:65:GLY:N	2.27	0.49
1:A:89:LEU:HD13	1:A:108:VAL:HG12	1.93	0.49
1:A:146:GLY:O	1:A:147:ARG:O	2.30	0.49
1:A:102:GLU:HG3	2:A:494:HOH:O	2.11	0.49
1:A:101:ASN:H	1:A:101:ASN:HD22	1.56	0.49
1:B:201:MET:CB	1:B:202:LYS:HD3	2.42	0.48
1:B:203:MET:O	1:B:206:GLN:OE1	2.32	0.48
1:B:226:THR:O	1:B:235:HIS:HA	2.14	0.48
1:B:229:SER:HG	1:B:295:ILE:HD12	1.77	0.47
1:B:266:PHE:CE2	1:B:270:LEU:HD12	2.50	0.47
1:B:313:TRP:CH2	1:B:334:LEU:HD12	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:254:LEU:HD22	1:B:332:TYR:CE1	2.49	0.47
1:B:240:GLU:OE2	1:B:307:ASP:HB3	2.14	0.47
1:B:223:LYS:HE2	2:B:477:HOH:O	2.13	0.47
1:A:119:ASP:O	1:A:119:ASP:OD1	2.33	0.47
1:B:247:VAL:HG12	1:B:251:ILE:HD12	1.96	0.47
1:A:36:TYR:CE1	1:A:107:ASP:HB3	2.50	0.47
1:B:296:LYS:NZ	1:B:296:LYS:HB2	2.23	0.47
1:B:299:LYS:H	1:B:299:LYS:HG3	1.60	0.47
1:A:88:GLU:O	1:A:88:GLU:HG3	2.15	0.47
1:B:203:MET:HE2	1:B:204:LEU:H	1.80	0.46
1:B:301:ASN:HB2	1:B:304:ILE:HB	1.98	0.46
1:A:54:LEU:HD23	1:A:139:ILE:HG21	1.96	0.46
1:B:215:PHE:CE1	1:B:221:LEU:HG	2.51	0.46
1:A:97:ARG:O	1:A:99:LYS:N	2.49	0.45
1:A:100:ILE:O	1:A:101:ASN:HB2	2.15	0.45
1:B:300:ILE:O	1:B:304:ILE:HG21	2.17	0.45
1:A:20:LEU:CD1	1:A:119:ASP:HB2	2.44	0.45
1:B:277:GLU:OE1	1:B:281:ASN:ND2	2.50	0.45
1:B:300:ILE:O	1:B:301:ASN:HB2	2.17	0.45
1:B:288:GLU:CB	1:B:291:LYS:HE2	2.47	0.45
1:B:299:LYS:HB2	1:B:299:LYS:HE3	1.49	0.45
1:B:279:PHE:CD2	1:B:283:LEU:HD11	2.52	0.45
1:B:275:ILE:HA	1:B:280:PHE:CE1	2.52	0.45
1:B:345:GLU:HA	1:B:345:GLU:OE1	2.16	0.45
1:B:288:GLU:HB2	1:B:291:LYS:HE2	1.99	0.44
1:B:291:LYS:HE3	1:B:292:PHE:CZ	2.51	0.44
1:A:3:MET:H	1:A:3:MET:HG3	1.40	0.44
1:A:122:ARG:O	1:B:328:ARG:NH2	2.51	0.44
1:B:205:GLU:HG2	1:B:221:LEU:HD13	1.99	0.44
1:B:203:MET:CE	1:B:204:LEU:H	2.31	0.44
1:B:333:LYS:HA	1:B:333:LYS:HD3	1.49	0.44
1:A:61:ASN:HD21	1:A:66:PHE:HB2	1.83	0.44
1:B:253:ASP:O	1:B:257:GLN:HB2	2.17	0.44
1:A:99:LYS:HG3	1:A:101:ASN:ND2	2.33	0.44
1:B:227:HIS:CE1	1:B:228:VAL:HG23	2.53	0.43
1:B:228:VAL:HG13	1:B:235:HIS:CG	2.53	0.43
1:A:54:LEU:HD21	1:A:136:LYS:HG3	2.00	0.43
1:B:228:VAL:HG13	1:B:235:HIS:NE2	2.34	0.43
1:A:135:PHE:O	1:A:139:ILE:HG13	2.19	0.43
1:B:274:LEU:HD13	1:B:339:ILE:HG23	2.01	0.42
1:B:285:GLN:NE2	2:B:646:HOH:O	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:211:LEU:HD21	1:B:312:LEU:HD21	2.01	0.42
1:A:2:LYS:HB2	1:A:2:LYS:HZ3	1.84	0.42
1:A:55:LEU:HD13	1:A:67:LEU:HD22	2.01	0.42
1:B:230:TYR:HB2	1:B:294:ARG:O	2.19	0.42
1:A:29:SER:HB2	1:A:97:ARG:HD2	2.01	0.42
1:A:27:HIS:HD2	1:A:36:TYR:OH	2.03	0.42
1:B:314:ALA:O	1:B:318:ILE:HG13	2.20	0.42
1:B:259:SER:HG	1:B:266:PHE:HE2	1.66	0.41
1:A:69:PRO:O	1:A:72:ALA:HB3	2.20	0.41
1:A:116:VAL:O	1:A:120:SER:HB3	2.20	0.41
1:B:258:TYR:CD2	1:B:340:LEU:HD21	2.56	0.41
1:A:93:ILE:CG1	1:A:108:VAL:HG21	2.39	0.41
1:B:261:ASN:ND2	1:B:266:PHE:HB2	2.30	0.41
1:B:309:PHE:O	1:B:313:TRP:HD1	2.04	0.41
1:B:296:LYS:O	1:B:297:ARG:HB2	2.20	0.41
1:A:86:LYS:HD3	2:A:487:HOH:O	2.21	0.41
1:B:313:TRP:CZ2	1:B:334:LEU:HD12	2.56	0.40
1:B:280:PHE:HA	1:B:283:LEU:HD12	2.02	0.40
1:B:243:GLY:HA2	1:B:246:LEU:HB2	2.04	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	145/147 (99%)	135 (93%)	8 (6%)	2 (1%)	14	7
1	B	145/147 (99%)	122 (84%)	20 (14%)	3 (2%)	9	3
All	All	290/294 (99%)	257 (89%)	28 (10%)	5 (2%)	11	4

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	98	GLY
1	A	101	ASN
1	B	231	SER
1	B	202	LYS
1	B	304	ILE

### 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	131/131 (100%)	109 (83%)	22 (17%)	2	1
1	B	131/131 (100%)	95 (72%)	36 (28%)	0	0
All	All	262/262 (100%)	204 (78%)	58 (22%)	1	0

All (58) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LYS
1	A	3	MET
1	A	9	LYS
1	A	16	LYS
1	A	18	LYS
1	A	21	LEU
1	A	42	LEU
1	A	46	LEU
1	A	51	ILE
1	A	55	LEU
1	A	61	ASN
1	A	62	LYS
1	A	64	GLU
1	A	70	LEU
1	A	91	LYS
1	A	97	ARG
1	A	100	ILE
1	A	101	ASN
1	A	103	THR

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Mol	Chain	Res	Type
1	A	130	LEU
1	A	137	GLU
1	A	144	LYS
1	B	201	MET
1	B	202	LYS
1	B	205	GLU
1	B	206	GLN
1	B	218	LYS
1	B	220	LEU
1	B	221	LEU
1	B	225	LEU
1	B	228	VAL
1	B	232	LYS
1	B	242	LEU
1	B	244	ASP
1	B	246	LEU
1	B	254	LEU
1	B	255	LEU
1	B	256	VAL
1	B	262	LYS
1	B	270	LEU
1	B	273	TYR
1	B	276	SER
1	B	282	LEU
1	B	296	LYS
1	B	297	ARG
1	B	299	LYS
1	B	300	ILE
1	B	304	ILE
1	B	305	ILE
1	B	307	ASP
1	B	318	ILE
1	B	322	ARG
1	B	330	LEU
1	B	333	LYS
1	B	341	SER
1	B	344	LYS
1	B	345	GLU
1	B	347	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	HIS
1	A	85	GLN
1	A	101	ASN
1	B	206	GLN
1	B	257	GLN
1	B	261	ASN
1	B	301	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](#)

There are no ligands in this entry.

### 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 ⓘ

EDS was not executed - this section will therefore be empty.

### 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.