



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 06:49 PM UTC

PDB ID : 3ORR / pdb_00003orr
Title : Crystal Structure of N5-Carboxyaminoimidazole synthetase from *Staphylococcus aureus*
Authors : Brugarolas, P.; Duguid, E.M.; Zhang, W.; Poor, C.B.; He, C.
Deposited on : 2010-09-07
Resolution : 2.23 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

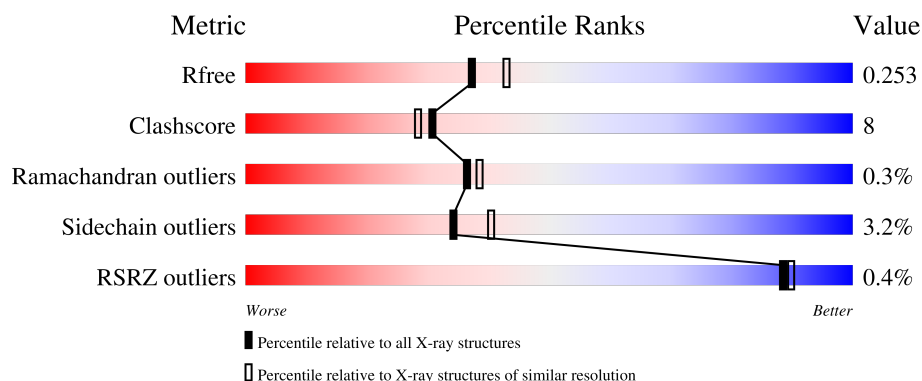
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.23 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	3416 (2.26-2.22)
Clashscore	190562	3556 (2.26-2.22)
Ramachandran outliers	187476	3500 (2.26-2.22)
Sidechain outliers	187428	3501 (2.26-2.22)
RSRZ outliers	180081	3415 (2.26-2.22)

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	377	% 80% 15% ..
1	B	377	79% 18% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6105 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 N5-carboxyaminoimidazole ribonucleotide synthetase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	368	Total	C	N	O	S	Se	0	0	0
			2945	1868	496	566	6	9			
1	B	367	Total	C	N	O	S	Se	0	1	0
			2945	1868	496	566	6	9			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP A6QFS4
A	-1	ASN	-	expression tag	UNP A6QFS4
A	0	ALA	-	expression tag	UNP A6QFS4
B	-2	SER	-	expression tag	UNP A6QFS4
B	-1	ASN	-	expression tag	UNP A6QFS4
B	0	ALA	-	expression tag	UNP A6QFS4

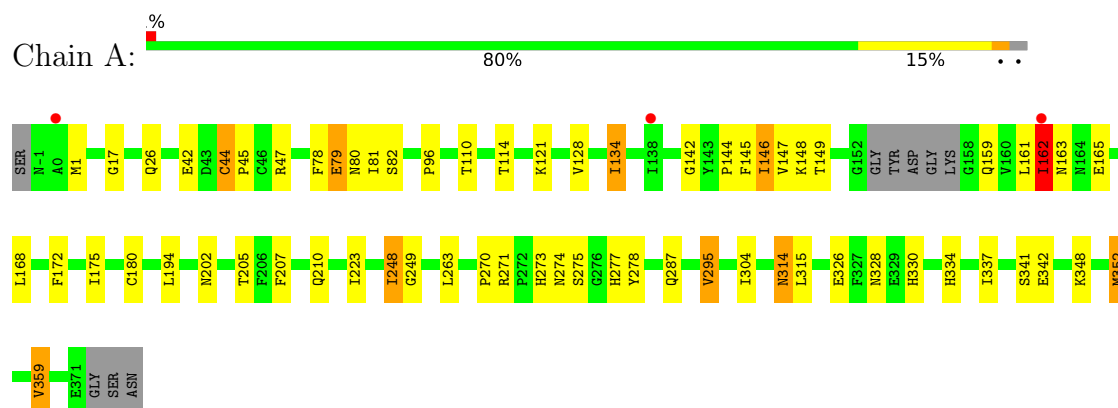
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	90	Total	O	0	0
			90	90		
2	B	125	Total	O	0	0
			125	125		

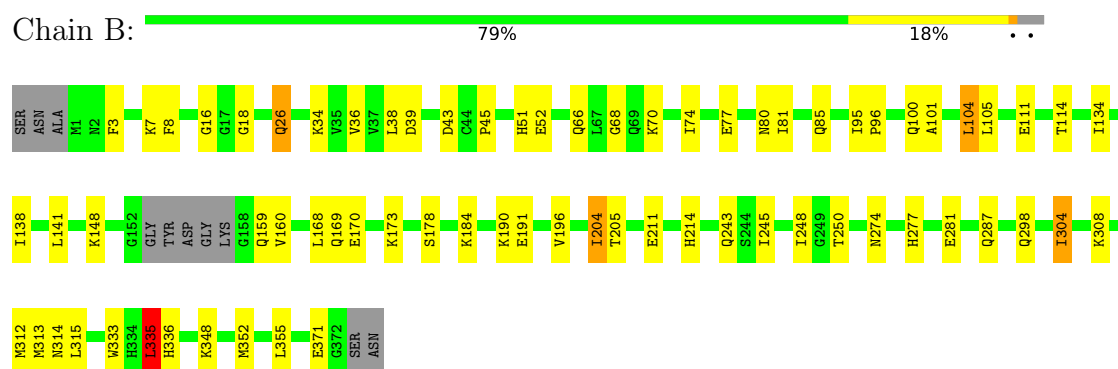
3 Residue-property plots

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: N5-carboxyaminoimidazole ribonucleotide synthetase



- Molecule 1: N5-carboxyaminoimidazole ribonucleotide synthetase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	66.11Å 53.33Å 107.41Å 90.00° 95.34° 90.00°	Depositor
Resolution (Å)	40.00 – 2.23 40.00 – 2.23	Depositor EDS
% Data completeness (in resolution range)	99.3 (40.00-2.23) 99.3 (40.00-2.23)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.27 (at 2.22Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.194 , 0.251 (Not available) , 0.253	Depositor DCC
R_{free} test set	1869 reflections (5.13%)	wwPDB-VP
Wilson B-factor (Å ²)	30.2	Xtriage
Anisotropy	0.107	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 14.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6105	wwPDB-VP
Average B, all atoms (Å ²)	11.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.67% 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

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	0.89	1/2989 (0.0%)	0.94	2/4016 (0.0%)
1	B	0.96	2/2990 (0.1%)	0.93	0/4018
All	All	0.93	3/5979 (0.1%)	0.94	2/8034 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	352	MSE	SE-CE	-5.97	1.77	1.95
1	B	335	LEU	C-O	-5.34	1.17	1.23
1	B	333	TRP	C-O	-5.02	1.17	1.24

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	44	CYS	CA-C-N	5.72	125.34	119.56
1	A	44	CYS	C-N-CA	5.72	125.34	119.56

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	2945	0	2930	52	0
1	B	2945	0	2929	48	0
2	A	90	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	125	0	0	5	0
All	All	6105	0	5859	97	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 8.

All (97) 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:315:LEU:HD21	1:B:352:MSE:HE3	1.27	1.13
1:A:26:GLN:HE21	1:B:26:GLN:NE2	1.73	0.86
1:A:26:GLN:HE21	1:B:26:GLN:HE21	1.29	0.81
1:A:78:PHE:HB3	1:A:80:ASN:ND2	2.02	0.74
1:A:146:ILE:HD11	1:A:159:GLN:HB3	1.74	0.70
1:B:101:ALA:O	1:B:105:LEU:HG	1.92	0.69
1:A:334:HIS:HD2	2:B:494:HOH:O	1.74	0.68
1:B:68:GLY:HA2	1:B:74:ILE:HD11	1.75	0.68
1:A:165:GLU:HA	1:A:168:LEU:HG	1.77	0.67
1:A:17:GLY:O	1:A:45:PRO:HD2	1.95	0.67
1:B:100:GLN:NE2	1:B:245:ILE:HA	2.12	0.64
1:B:336:HIS:HD2	2:B:453:HOH:O	1.78	0.64
1:B:308:LYS:HD3	1:B:355:LEU:HB3	1.80	0.64
1:B:315:LEU:HD21	1:B:352:MSE:CE	2.18	0.63
1:B:138:ILE:HD11	1:B:168:LEU:HD21	1.79	0.63
1:B:274:ASN:HD22	1:B:348:LYS:NZ	1.94	0.63
1:B:214:HIS:HE1	2:B:474:HOH:O	1.82	0.62
1:A:273:HIS:HD2	1:A:275:SER:OG	1.85	0.60
1:B:104:LEU:HD22	1:B:111:GLU:HA	1.84	0.60
1:A:148:LYS:HD3	1:A:159:GLN:HE21	1.66	0.59
1:A:271:ARG:HB2	2:A:451:HOH:O	2.01	0.59
1:B:134:ILE:O	1:B:138:ILE:HG12	2.01	0.59
1:A:274:ASN:HD22	1:A:348:LYS:NZ	2.02	0.58
1:B:315:LEU:CD2	1:B:352:MSE:HE3	2.18	0.58
1:A:78:PHE:HB3	1:A:80:ASN:HD21	1.68	0.57
1:A:271:ARG:HG3	1:A:271:ARG:HH11	1.70	0.56
1:B:205:THR:HG22	1:B:304:ILE:HD12	1.87	0.56
1:B:277:HIS:HE1	2:B:454:HOH:O	1.88	0.56
1:A:96:PRO:HB2	1:A:248:ILE:HG12	1.87	0.55
1:B:281:GLU:OE2	1:B:336:HIS:HE1	1.89	0.55
1:B:274:ASN:HD21	1:B:314:ASN:HD21	1.53	0.55
1:A:248:ILE:HD12	1:A:249:GLY:N	2.21	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:334:HIS:CD2	2:B:494:HOH:O	2.53	0.54
1:A:134:ILE:HG21	1:A:172:PHE:HE1	1.72	0.54
1:B:196:VAL:HG11	1:B:204:ILE:HD11	1.89	0.54
1:B:16:GLY:O	1:B:77:GLU:HG2	2.07	0.53
1:B:81:ILE:HD11	1:B:85:GLN:HG3	1.90	0.53
1:B:274:ASN:ND2	1:B:314:ASN:HD21	2.07	0.53
1:B:52:GLU:OE2	1:B:70:LYS:HE2	2.11	0.51
1:A:147:VAL:HG11	1:A:175:ILE:HD11	1.93	0.51
1:B:66:GLN:O	1:B:70:LYS:HG3	2.11	0.51
1:A:207:PHE:CD1	1:A:304:ILE:HD13	2.46	0.51
1:A:273:HIS:H	1:A:287:GLN:HE22	1.59	0.50
1:A:121:LYS:HD2	1:A:263:LEU:O	2.11	0.50
1:B:141:LEU:O	1:B:184:LYS:HE2	2.12	0.50
1:A:273:HIS:CD2	1:A:275:SER:H	2.31	0.49
1:B:335:LEU:CD1	1:B:352:MSE:HE2	2.42	0.49
1:A:274:ASN:HD22	1:A:348:LYS:HZ2	1.61	0.49
1:A:110:THR:O	1:A:114:THR:HG23	2.12	0.49
1:B:3:PHE:CE2	1:B:248:ILE:HG13	2.47	0.48
1:B:38:LEU:C	1:B:38:LEU:HD23	2.38	0.48
1:B:308:LYS:CD	1:B:355:LEU:HB3	2.43	0.48
1:B:190:LYS:HE3	1:B:211:GLU:OE2	2.14	0.48
1:A:144:PRO:HB3	1:A:163:ASN:HD22	1.79	0.48
1:A:326:GLU:O	1:A:330:HIS:HD2	1.97	0.48
1:B:18:GLY:HA2	1:B:45:PRO:HG2	1.96	0.48
1:B:52:GLU:CD	1:B:70:LYS:HE2	2.39	0.47
1:A:194:LEU:HD12	2:A:427:HOH:O	2.14	0.47
1:A:277:HIS:HE1	2:A:411:HOH:O	1.97	0.47
1:A:145:PHE:CE1	1:A:162:ILE:HD12	2.51	0.46
1:A:128:VAL:HG11	1:A:134:ILE:HD12	1.98	0.46
1:A:1:MSE:HE1	1:A:202:ASN:HB2	1.98	0.45
1:A:314:ASN:HD22	1:A:314:ASN:HA	1.63	0.45
1:B:36:VAL:HG22	1:B:52:GLU:HB3	1.97	0.45
1:A:148:LYS:HG2	1:A:159:GLN:HG2	1.99	0.45
1:A:81:ILE:HG12	1:A:82:SER:O	2.17	0.45
1:B:148:LYS:HD3	1:B:159:GLN:HE21	1.82	0.45
1:A:128:VAL:HG12	1:A:175:ILE:HG12	1.99	0.44
1:A:223:ILE:HG21	1:A:359:VAL:HG11	1.99	0.44
1:A:248:ILE:HD11	1:A:295:VAL:HA	1.98	0.44
1:A:315:LEU:HD11	1:A:352:MSE:HE3	2.00	0.44
1:B:74:ILE:HB	1:B:95:ILE:HG12	2.00	0.44
1:B:191:GLU:OE1	1:B:214:HIS:HD2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:104:LEU:HD11	1:B:114:THR:HG21	1.99	0.44
1:A:142:GLY:HA3	1:A:145:PHE:HD2	1.83	0.44
1:A:147:VAL:HG12	1:A:180:CYS:HB3	1.99	0.43
1:B:160:VAL:HG21	1:B:170:GLU:HG2	2.00	0.43
1:B:169:GLN:HG2	1:B:173:LYS:HE3	2.00	0.43
1:B:312:MSE:HG2	1:B:313:MSE:N	2.32	0.43
1:A:205:THR:HG22	1:A:304:ILE:HD12	2.01	0.43
1:B:7:LYS:O	1:B:8:PHE:C	2.61	0.43
1:B:250:THR:HG21	1:B:287:GLN:HB2	2.01	0.43
1:A:44:CYS:HA	1:A:45:PRO:HD3	1.84	0.42
1:A:1:MSE:SE	1:A:202:ASN:HB2	2.69	0.42
1:B:196:VAL:CG1	1:B:204:ILE:HD11	2.50	0.42
1:A:79:GLU:OE2	1:A:270:PRO:HD2	2.19	0.42
1:A:223:ILE:CG2	1:A:359:VAL:HG11	2.49	0.41
1:A:145:PHE:O	1:A:162:ILE:HB	2.21	0.41
1:B:43:ASP:CG	1:B:43:ASP:O	2.63	0.41
1:A:17:GLY:O	1:A:44:CYS:HB2	2.20	0.41
1:A:328:ASN:HA	1:B:298:GLN:NE2	2.36	0.41
1:B:38:LEU:HD23	1:B:39:ASP:N	2.34	0.41
1:A:1:MSE:CE	1:A:202:ASN:HB2	2.50	0.41
1:B:34:LYS:HA	1:B:51:HIS:CE1	2.55	0.41
1:A:42:GLU:HG2	1:A:47:ARG:HH12	1.86	0.40
1:A:146:ILE:HD12	1:A:161:LEU:HA	2.02	0.40
1:A:210:GLN:HG2	1:A:278:TYR:CE1	2.57	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	364/377 (97%)	352 (97%)	10 (3%)	2 (0%)	24 23

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	364/377 (97%)	350 (96%)	14 (4%)	0	100	100
All	All	728/754 (97%)	702 (96%)	24 (3%)	2 (0%)	36	38

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	162	ILE
1	A	79	GLU

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	327/324 (101%)	316 (97%)	11 (3%)	32	38
1	B	327/324 (101%)	317 (97%)	10 (3%)	35	41
All	All	654/648 (101%)	633 (97%)	21 (3%)	34	40

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	134	ILE
1	A	146	ILE
1	A	149	THR
1	A	162	ILE
1	A	248	ILE
1	A	295	VAL
1	A	314	ASN
1	A	337	ILE
1	A	341	SER
1	A	342	GLU
1	A	359	VAL
1	B	26	GLN
1	B	80	ASN
1	B	96	PRO

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Mol	Chain	Res	Type
1	B	104	LEU
1	B	178	SER
1	B	204	ILE
1	B	243	GLN
1	B	304	ILE
1	B	335	LEU
1	B	371	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	2	ASN
1	A	80	ASN
1	A	84	GLN
1	A	159	GLN
1	A	210	GLN
1	A	212	ASN
1	A	273	HIS
1	A	274	ASN
1	A	277	HIS
1	A	287	GLN
1	A	302	ASN
1	A	314	ASN
1	A	330	HIS
1	A	334	HIS
1	B	26	GLN
1	B	29	GLN
1	B	100	GLN
1	B	159	GLN
1	B	214	HIS
1	B	217	GLN
1	B	243	GLN
1	B	262	GLN
1	B	274	ASN
1	B	277	HIS
1	B	334	HIS
1	B	336	HIS

5.3.3 RNA ⓘ

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](#)

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 [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	359/377 (95%)	-0.30	3 (0%) 82 84	2, 6, 45, 54	0
1	B	358/377 (94%)	-0.46	0 100 100	2, 3, 40, 49	1 (0%)
All	All	717/754 (95%)	-0.38	3 (0%) 88 90	2, 4, 43, 54	1 (0%)

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	0	ALA	2.5
1	A	138	ILE	2.1
1	A	162	ILE	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.