



wwPDB X-ray Structure Validation Summary Report ⓘ

May 23, 2020 – 03:49 am BST

PDB ID : 5ZLI
Title : Crystal structure of glutamine synthetase from helicobacter pylori
Authors : Joo, H.K.; Lee, J.Y.
Deposited on : 2018-03-28
Resolution : 2.80 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

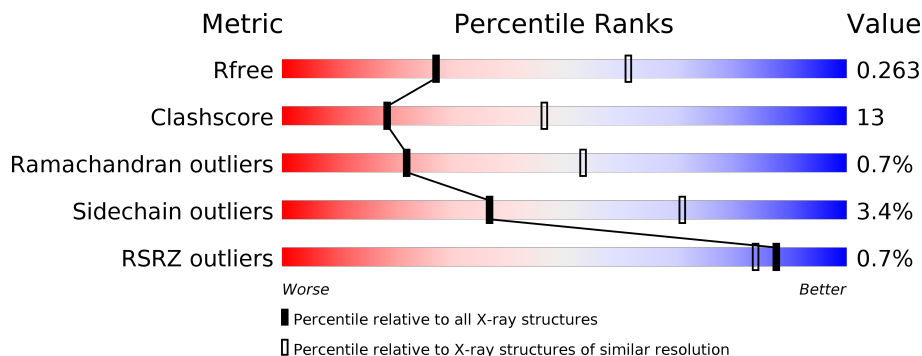
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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 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	481	
1	B	481	
1	C	481	
1	D	481	
1	E	481	
1	F	481	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 22534 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.

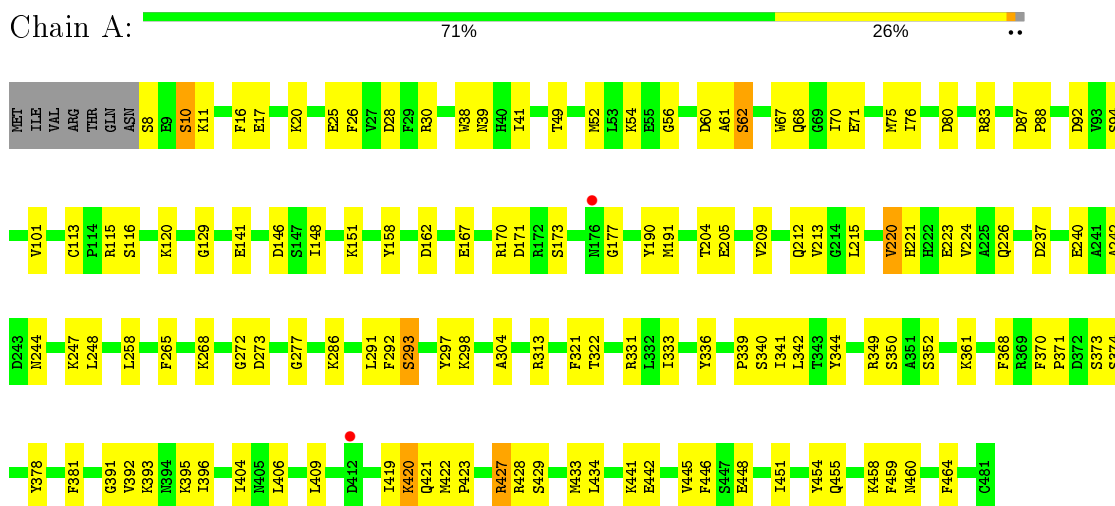
- Molecule 1 is a protein called Glutamine synthetase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	474	3752	2404	625	705	18	0	0	0
1	B	475	3760	2408	627	707	18	0	0	0
1	C	474	3752	2404	625	705	18	0	0	0
1	D	475	3760	2408	627	707	18	0	0	0
1	E	474	3758	2407	628	705	18	0	0	0
1	F	474	3752	2404	625	705	18	0	0	0

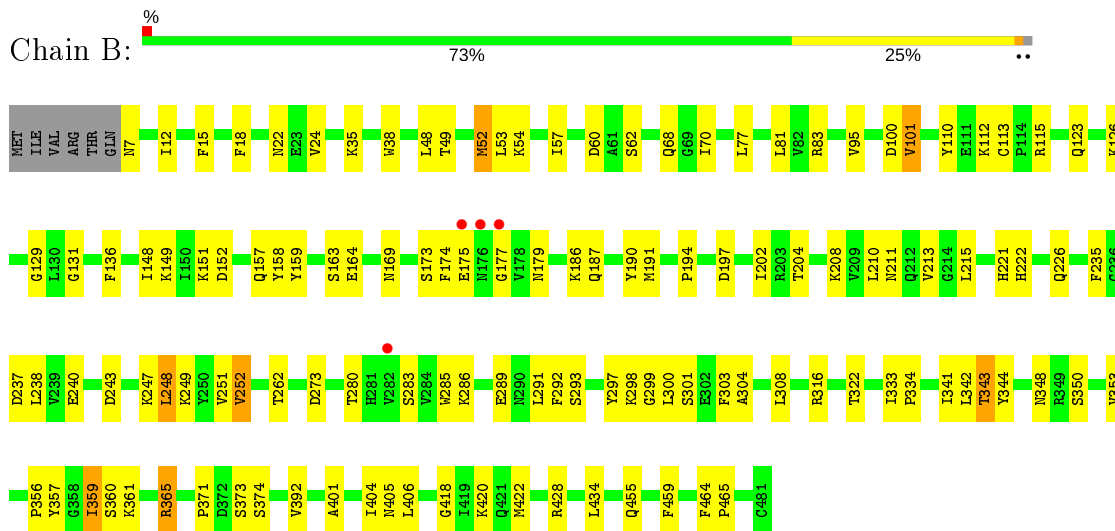
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: Glutamine synthetase

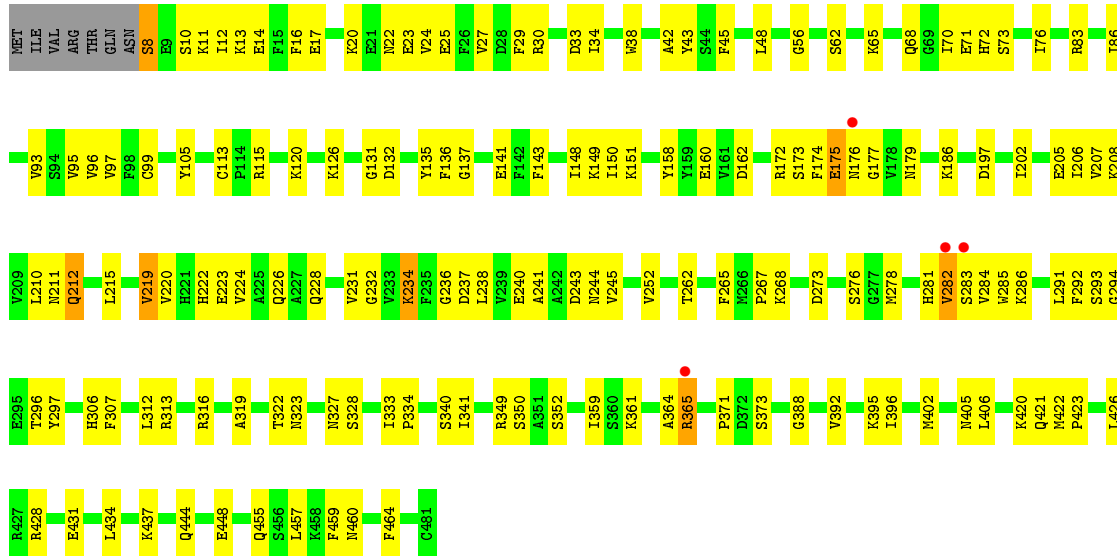


- Molecule 1: Glutamine synthetase

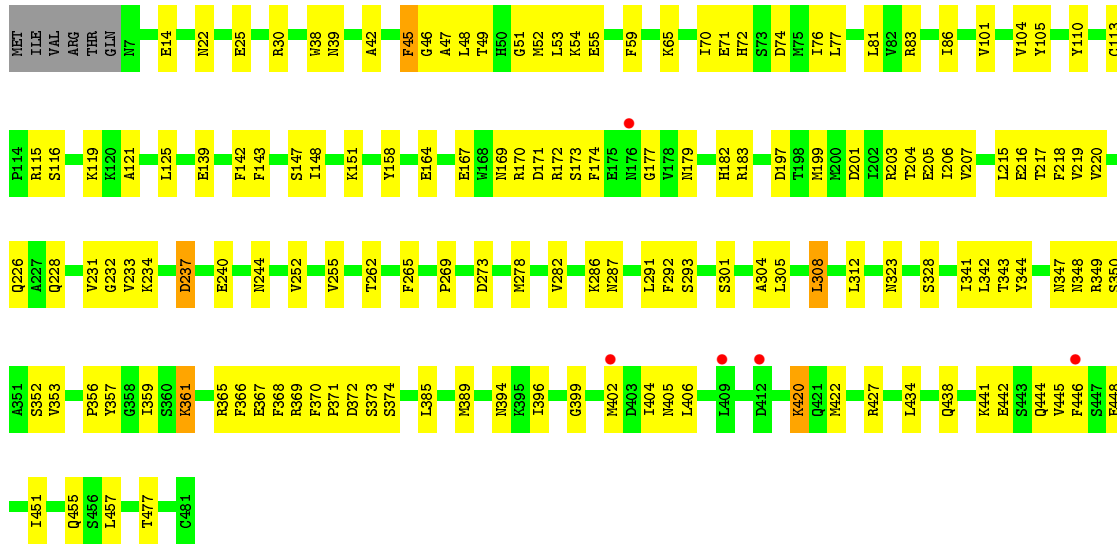


- Molecule 1: Glutamine synthetase

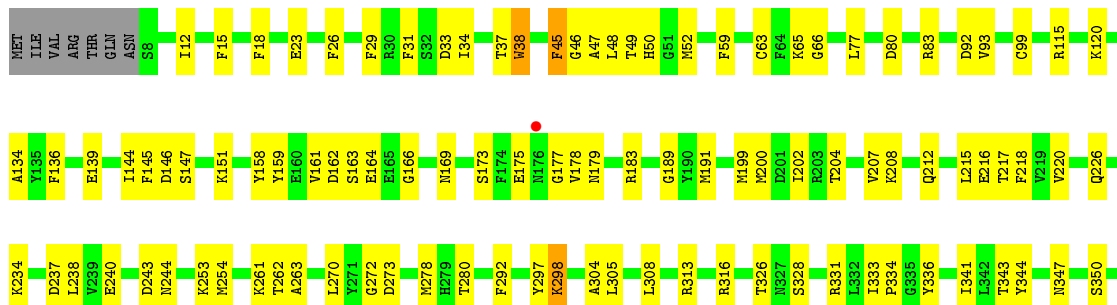


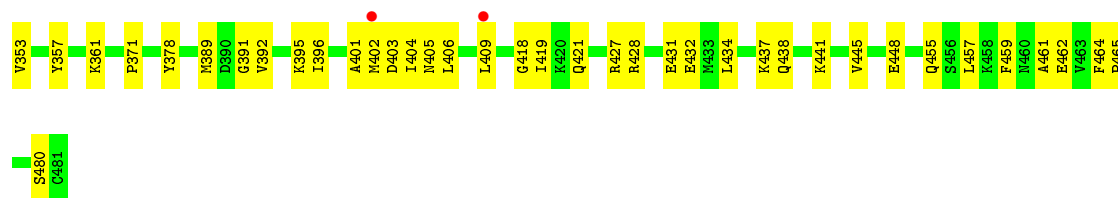


• Molecule 1: Glutamine synthetase

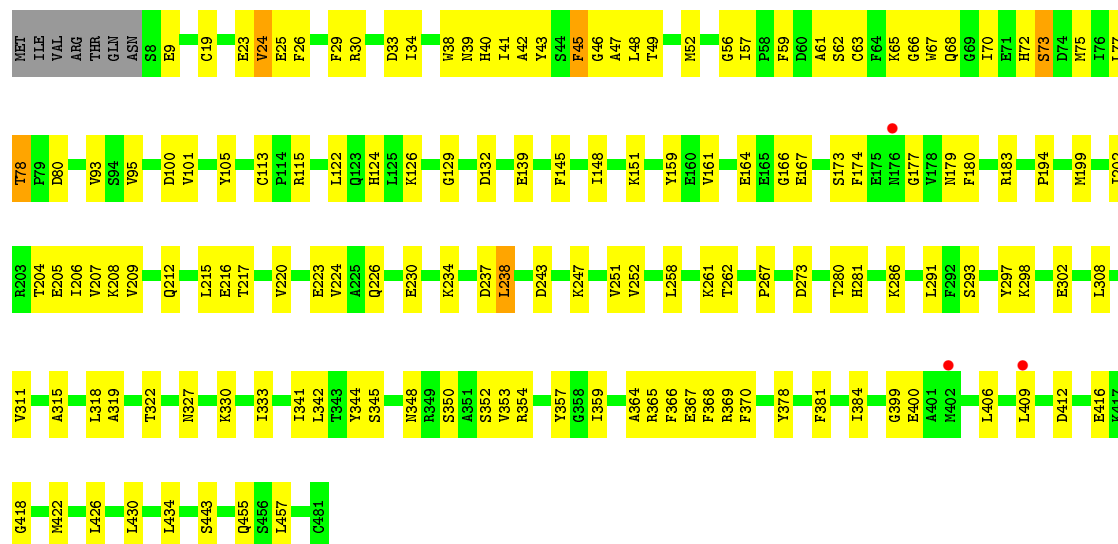


• Molecule 1: Glutamine synthetase





● Molecule 1: Glutamine synthetase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	191.09Å 131.46Å 131.41Å 90.00° 122.67° 90.00°	Depositor
Resolution (Å)	20.00 – 2.80 46.76 – 2.79	Depositor EDS
% Data completeness (in resolution range)	99.6 (20.00-2.80) 99.2 (46.76-2.79)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.42 (at 2.81Å)	Xtrriage
Refinement program	PHENIX 1.10.1_2155	Depositor
R, R_{free}	0.190 , 0.260 0.193 , 0.263	Depositor DCC
R_{free} test set	2000 reflections (2.96%)	wwPDB-VP
Wilson B-factor (Å ²)	49.2	Xtrriage
Anisotropy	0.080	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 40.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.39$, $\langle L^2 \rangle = 0.22$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	22534	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.27% 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](#)

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	0.56	0/3852	0.74	3/5207 (0.1%)
1	B	0.55	0/3860	0.70	1/5218 (0.0%)
1	C	0.59	2/3852 (0.1%)	0.73	2/5207 (0.0%)
1	D	0.53	0/3860	0.70	1/5218 (0.0%)
1	E	0.55	1/3858 (0.0%)	0.70	0/5214
1	F	0.58	0/3852	0.74	1/5207 (0.0%)
All	All	0.56	3/23134 (0.0%)	0.72	8/31271 (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	C	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	45	PHE	CE1-CZ	6.59	1.49	1.37
1	E	63	CYS	CB-SG	-5.74	1.72	1.81
1	C	113	CYS	CB-SG	5.14	1.91	1.82

The worst 5 of 8 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	258	LEU	CA-CB-CG	6.97	131.33	115.30
1	F	238	LEU	CB-CG-CD1	-6.87	99.33	111.00
1	B	248	LEU	CA-CB-CG	5.87	128.81	115.30
1	A	427	ARG	NE-CZ-NH1	-5.77	117.41	120.30
1	D	305	LEU	CA-CB-CG	5.54	128.04	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	174	PHE	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	3752	0	3580	98	0
1	B	3760	0	3586	78	0
1	C	3752	0	3580	130	0
1	D	3760	0	3586	113	0
1	E	3758	0	3591	93	0
1	F	3752	0	3580	108	0
All	All	22534	0	21503	581	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 13.

The worst 5 of 581 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:C:293:SER:HB2	1:C:361:LYS:HD2	1.42	1.01
1:F:238:LEU:HD11	1:F:381:PHE:HB3	1.41	0.98
1:F:122:LEU:HD21	1:F:238:LEU:HB3	1.45	0.97
1:A:167:GLU:OE1	1:A:170:ARG:NH1	2.02	0.93
1:C:283:SER:HA	1:C:365:ARG:CZ	2.04	0.87

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	472/481 (98%)	443 (94%)	26 (6%)	3 (1%)	25	56
1	B	473/481 (98%)	447 (94%)	23 (5%)	3 (1%)	25	56
1	C	472/481 (98%)	442 (94%)	28 (6%)	2 (0%)	34	66
1	D	473/481 (98%)	447 (94%)	23 (5%)	3 (1%)	25	56
1	E	472/481 (98%)	440 (93%)	27 (6%)	5 (1%)	14	41
1	F	472/481 (98%)	440 (93%)	27 (6%)	5 (1%)	14	41
All	All	2834/2886 (98%)	2659 (94%)	154 (5%)	21 (1%)	22	53

5 of 21 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	298	LYS
1	D	45	PHE
1	D	46	GLY
1	E	45	PHE
1	E	46	GLY

5.3.2 Protein sidechains [i](#)

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	398/415 (96%)	386 (97%)	12 (3%)	41	75
1	B	399/415 (96%)	382 (96%)	17 (4%)	29	62
1	C	398/415 (96%)	383 (96%)	15 (4%)	33	67
1	D	399/415 (96%)	386 (97%)	13 (3%)	38	72
1	E	399/415 (96%)	391 (98%)	8 (2%)	55	84
1	F	398/415 (96%)	381 (96%)	17 (4%)	29	62
All	All	2391/2490 (96%)	2309 (97%)	82 (3%)	37	71

5 of 82 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	237	ASP
1	D	38	TRP
1	F	216	GLU
1	C	296	THR
1	C	365	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	124	HIS
1	D	287	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 [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	474/481 (98%)	-0.42	2 (0%) 92 91	21, 39, 63, 99	0
1	B	475/481 (98%)	-0.35	4 (0%) 86 81	22, 41, 60, 77	0
1	C	474/481 (98%)	-0.30	4 (0%) 86 81	24, 43, 72, 86	0
1	D	475/481 (98%)	-0.23	5 (1%) 80 75	26, 48, 72, 97	0
1	E	474/481 (98%)	-0.35	3 (0%) 89 86	24, 43, 66, 84	0
1	F	474/481 (98%)	-0.38	3 (0%) 89 86	22, 39, 63, 104	0
All	All	2846/2886 (98%)	-0.34	21 (0%) 87 84	21, 42, 67, 104	0

The worst 5 of 21 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	409	LEU	4.5
1	C	283	SER	3.5
1	D	176	ASN	3.0
1	F	176	ASN	3.0
1	D	446	PHE	2.9

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](#)

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.