



Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 05:12 am BST

PDB ID : 4K86
Title : Crystal structure of human prolyl-tRNA synthetase (apo form)
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Deposited on : 2013-04-18
Resolution : 2.40 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
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

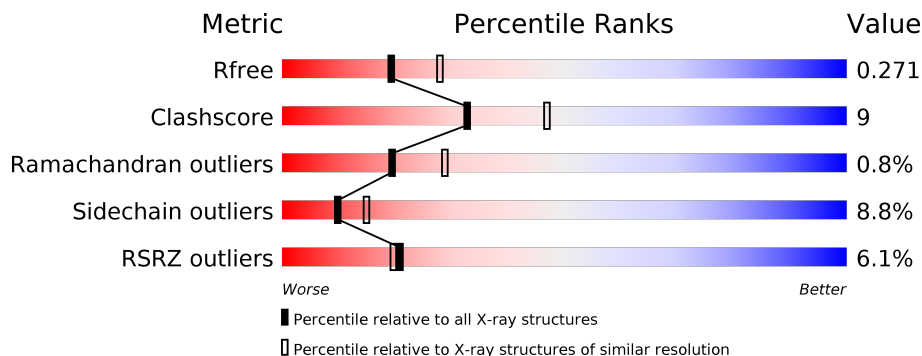
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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	535	 6% (red), 67% (green), 21% (yellow), 9% (orange), 9% (grey)

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 3922 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 Proline-tRNA ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	488	3914	2505	662	722	25	0	0	0

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-22	MET	-	expression tag	UNP P07814
A	-21	GLY	-	expression tag	UNP P07814
A	-20	SER	-	expression tag	UNP P07814
A	-19	SER	-	expression tag	UNP P07814
A	-18	HIS	-	expression tag	UNP P07814
A	-17	HIS	-	expression tag	UNP P07814
A	-16	HIS	-	expression tag	UNP P07814
A	-15	HIS	-	expression tag	UNP P07814
A	-14	HIS	-	expression tag	UNP P07814
A	-13	HIS	-	expression tag	UNP P07814
A	-12	SER	-	expression tag	UNP P07814
A	-11	SER	-	expression tag	UNP P07814
A	-10	GLY	-	expression tag	UNP P07814
A	-9	LEU	-	expression tag	UNP P07814
A	-8	VAL	-	expression tag	UNP P07814
A	-7	PRO	-	expression tag	UNP P07814
A	-6	ARG	-	expression tag	UNP P07814
A	-5	GLY	-	expression tag	UNP P07814
A	-4	SER	-	expression tag	UNP P07814
A	-3	HIS	-	expression tag	UNP P07814
A	-2	MET	-	expression tag	UNP P07814
A	-1	ALA	-	expression tag	UNP P07814

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total 1	Zn 1	0	0

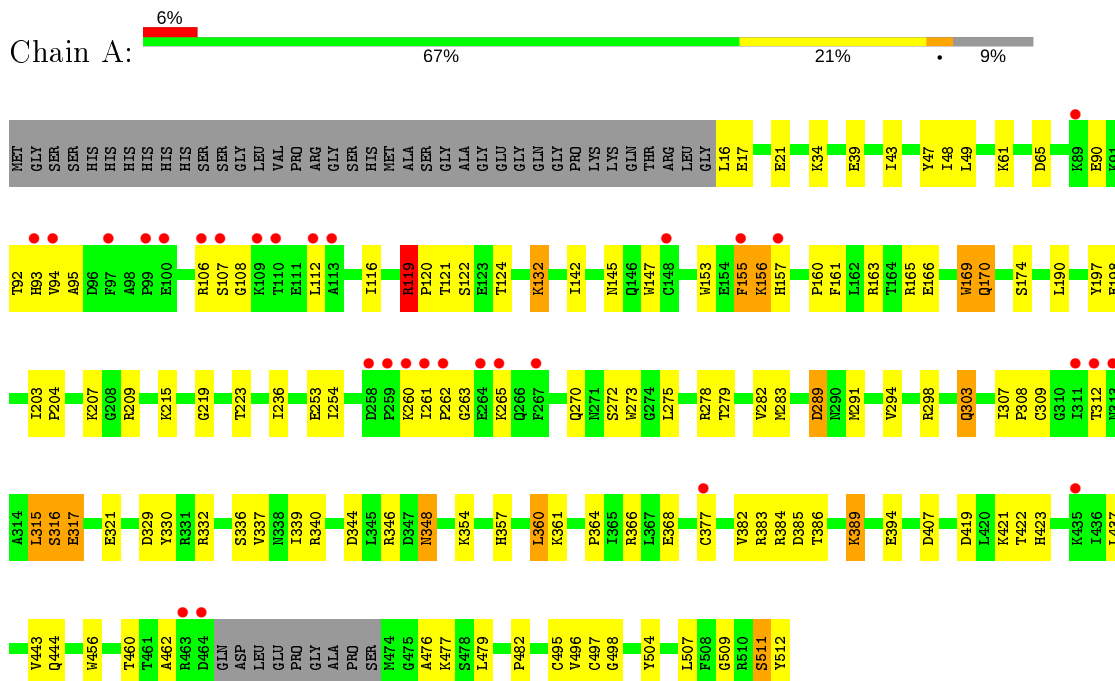
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	7	Total 7	O 7	0	0

3 Residue-property plots [i](#)

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: Proline-tRNA ligase



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	119.85Å 119.85Å 95.72Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	39.23 – 2.40 39.23 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.7 (39.23-2.40) 99.6 (39.23-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.83 (at 2.29Å)	Xtrriage
Refinement program	PHENIX 1.7.3_928	Depositor
R, R_{free}	0.239 , 0.277 0.231 , 0.271	Depositor DCC
R_{free} test set	2009 reflections (5.67%)	wwPDB-VP
Wilson B-factor (Å ²)	33.6	Xtrriage
Anisotropy	0.297	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 66.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.39$, $\langle L^2 \rangle = 0.22$	Xtrriage
Estimated twinning fraction	0.064 for -h,-k,l	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3922	wwPDB-VP
Average B, all atoms (Å ²)	65.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.36% 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:
ZN

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	0.33	1/4007 (0.0%)	0.52	1/5425 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	377	CYS	CB-SG	5.41	1.91	1.82

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	119	ARG	C-N-CD	-6.37	106.58	120.60

There are no chirality outliers.

There are no planarity outliers.

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	3914	0	3879	68	0
2	A	1	0	0	0	0
3	A	7	0	0	2	0
All	All	3922	0	3879	68	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 9.

All (68) 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:34:LYS:HB3	1:A:360:LEU:HD11	1.69	0.73
1:A:145:ASN:OD1	1:A:170:GLN:NE2	2.22	0.71
1:A:348:ASN:OD1	1:A:348:ASN:N	2.20	0.67
1:A:329:ASP:OD1	1:A:332:ARG:NH2	2.29	0.65
1:A:106:ARG:NH1	1:A:108:GLY:O	2.29	0.65
1:A:337:VAL:HG23	1:A:339:ILE:HG13	1.81	0.62
1:A:346:ARG:H	1:A:354:LYS:HZ1	1.45	0.61
1:A:479:LEU:HD12	1:A:507:LEU:HG	1.82	0.60
1:A:215:LYS:NZ	1:A:219:GLY:O	2.38	0.57
1:A:203:ILE:HD11	1:A:283:MET:HA	1.88	0.56
1:A:261:ILE:HG22	1:A:263:GLY:HA2	1.87	0.56
1:A:253:GLU:OE1	1:A:265:LYS:NZ	2.33	0.55
1:A:107:SER:HB3	1:A:112:LEU:HD21	1.88	0.55
1:A:289:ASP:HB3	1:A:291:MET:H	1.72	0.54
1:A:190:LEU:HD22	1:A:223:THR:HB	1.89	0.54
1:A:346:ARG:H	1:A:354:LYS:NZ	2.05	0.53
1:A:383:ARG:NH2	1:A:407:ASP:OD2	2.42	0.53
1:A:92:THR:OG1	1:A:93:HIS:N	2.41	0.53
1:A:511:SER:OG	1:A:512:TYR:N	2.39	0.52
1:A:203:ILE:HD13	1:A:282:VAL:HG12	1.91	0.52
1:A:344:ASP:O	1:A:354:LYS:NZ	2.38	0.52
1:A:39:GLU:HB3	1:A:48:ILE:HD12	1.90	0.52
1:A:155:PHE:HD1	1:A:165:ARG:HH22	1.56	0.51
1:A:456:TRP:O	1:A:460:THR:HG22	2.11	0.51
1:A:497:CYS:SG	1:A:498:GLY:N	2.83	0.51
1:A:315:LEU:HD22	1:A:316:SER:H	1.76	0.50
1:A:147:TRP:CE3	1:A:170:GLN:HB3	2.46	0.49
1:A:385:ASP:OD1	1:A:386:THR:N	2.39	0.49
1:A:197:TYR:CZ	1:A:279:THR:HG22	2.48	0.47
1:A:289:ASP:HB2	1:A:384:ARG:HH22	1.79	0.47
1:A:303:GLN:HA	1:A:364:PRO:HD2	1.95	0.47
1:A:495:CYS:SG	1:A:496:VAL:N	2.88	0.47
1:A:61:LYS:O	1:A:65:ASP:HB2	2.15	0.47
1:A:90:GLU:OE2	1:A:93:HIS:ND1	2.28	0.46
1:A:477:LYS:HG2	3:A:805:HOH:O	2.16	0.46
1:A:236:ILE:HD11	1:A:278:ARG:HD3	1.98	0.45
1:A:142:ILE:O	1:A:174:SER:HA	2.17	0.45
1:A:21:GLU:OE1	1:A:21:GLU:N	2.48	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:444:GLN:NE2	1:A:482:PRO:HG3	2.31	0.45
1:A:49:LEU:HD23	1:A:49:LEU:HA	1.74	0.45
1:A:437:LEU:HD11	1:A:509:GLY:HA2	1.98	0.44
1:A:107:SER:OG	1:A:107:SER:O	2.35	0.44
1:A:161:PHE:N	3:A:804:HOH:O	2.46	0.44
1:A:169:TRP:HB2	1:A:275:LEU:O	2.18	0.44
1:A:357:HIS:O	1:A:361:LYS:HG3	2.18	0.43
1:A:94:VAL:HG12	1:A:95:ALA:O	2.18	0.43
1:A:254:ILE:HD13	1:A:270:GLN:OE1	2.19	0.43
1:A:422:THR:HG22	1:A:423:HIS:ND1	2.34	0.43
1:A:366:ARG:NH1	1:A:382:VAL:HG21	2.33	0.43
1:A:389:LYS:HB3	1:A:389:LYS:HE2	1.68	0.43
1:A:303:GLN:HG2	1:A:340:ARG:HB2	2.01	0.42
1:A:34:LYS:O	1:A:360:LEU:HD21	2.19	0.42
1:A:47:TYR:CZ	1:A:160:PRO:HG3	2.55	0.42
1:A:198:GLU:HG2	1:A:204:PRO:HA	2.01	0.42
1:A:316:SER:OG	1:A:317:GLU:N	2.53	0.42
1:A:170:GLN:HB2	1:A:170:GLN:HE21	1.71	0.42
1:A:132:LYS:HB3	1:A:132:LYS:HE2	1.89	0.41
1:A:330:TYR:CE1	1:A:394:GLU:HG3	2.55	0.41
1:A:47:TYR:CE2	1:A:160:PRO:HG3	2.55	0.41
1:A:383:ARG:HH21	1:A:407:ASP:HB3	1.86	0.41
1:A:121:THR:OG1	1:A:124:THR:OG1	2.27	0.41
1:A:190:LEU:HD21	1:A:207:LYS:HG2	2.03	0.41
1:A:444:GLN:HE21	1:A:482:PRO:HG3	1.86	0.41
1:A:366:ARG:NH1	1:A:368:GLU:OE2	2.54	0.41
1:A:366:ARG:HB3	1:A:382:VAL:HB	2.03	0.41
1:A:437:LEU:HD21	1:A:476:ALA:HB2	2.03	0.41
1:A:156:LYS:HB2	1:A:163:ARG:NH1	2.36	0.40
1:A:307:ILE:HA	1:A:308:PRO:HD2	1.81	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	484/535 (90%)	455 (94%)	25 (5%)	4 (1%)	19 29

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	119	ARG
1	A	262	PRO
1	A	462	ALA
1	A	120	PRO

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	421/456 (92%)	384 (91%)	37 (9%)	10 15

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	LEU
1	A	17	GLU
1	A	43	ILE
1	A	116	ILE
1	A	119	ARG
1	A	122	SER
1	A	132	LYS
1	A	153	TRP
1	A	155	PHE
1	A	156	LYS
1	A	157	HIS
1	A	166	GLU
1	A	169	TRP
1	A	170	GLN
1	A	209	ARG

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Mol	Chain	Res	Type
1	A	260	LYS
1	A	272	SER
1	A	273	TRP
1	A	289	ASP
1	A	294	VAL
1	A	298	ARG
1	A	303	GLN
1	A	309	CYS
1	A	312	THR
1	A	315	LEU
1	A	316	SER
1	A	317	GLU
1	A	321	GLU
1	A	336	SER
1	A	348	ASN
1	A	360	LEU
1	A	389	LYS
1	A	419	ASP
1	A	421	LYS
1	A	443	VAL
1	A	504	TYR
1	A	511	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	173	HIS

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

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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, 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	488/535 (91%)	0.09	30 (6%) 21 20	28, 57, 123, 147	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	260	LYS	6.0
1	A	264	GLU	6.0
1	A	261	ILE	5.7
1	A	106	ARG	5.6
1	A	113	ALA	5.4
1	A	259	PRO	5.2
1	A	109	LYS	5.0
1	A	313	ASN	4.8
1	A	157	HIS	4.8
1	A	312	THR	4.7
1	A	463	ARG	4.6
1	A	267	PHE	3.9
1	A	97	PHE	3.9
1	A	311	ILE	3.8
1	A	464	ASP	3.8
1	A	265	LYS	3.8
1	A	93	HIS	3.2
1	A	89	LYS	3.2
1	A	100	GLU	3.1
1	A	112	LEU	2.8
1	A	110	THR	2.6
1	A	262	PRO	2.6
1	A	99	PRO	2.6
1	A	377	CYS	2.4
1	A	435	LYS	2.4
1	A	258	ASP	2.3
1	A	148	CYS	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	155	PHE	2.2
1	A	107	SER	2.1
1	A	94	VAL	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. 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
2	ZN	A	700	1/1	0.92	0.27	77,77,77,77	0

6.5 Other polymers [i](#)

There are no such residues in this entry.