



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 4, 2016 – 10:55 AM EDT

PDB ID : 4KQE
Title : The mutant structure of the human glycyl-tRNA synthetase E71G
Authors : Qin, X.; Hao, Z.; Tian, Q.; Zhang, Z.; Zhou, C.; Xie, W.
Deposited on : 2013-05-15
Resolution : 2.74 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
<http://wwpdb.org/validation/2016/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
Mogul : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027107
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : rb-20027107

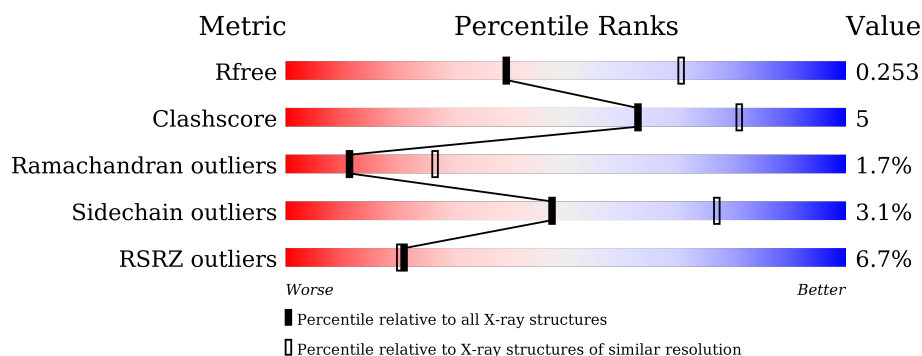
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.74 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}	91344	3050 (2.78-2.70)
Clashscore	102246	3424 (2.78-2.70)
Ramachandran outliers	100387	3367 (2.78-2.70)
Sidechain outliers	100360	3368 (2.78-2.70)
RSRZ outliers	91569	3055 (2.78-2.70)

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. 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	693	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	GOL	A	701	-	-	-	X

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 4756 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.

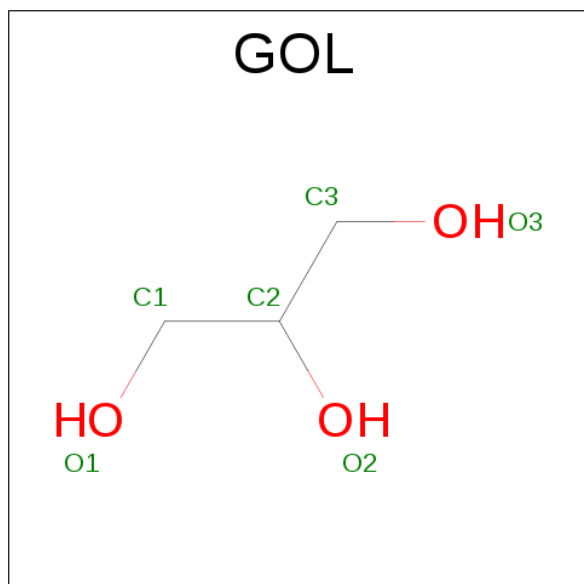
- Molecule 1 is a protein called Glycine-tRNA ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	597	4650	2953	793	875	29	0	0	0

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	71	GLY	GLU	ENGINEERED MUTATION	UNP P41250
A	686	LEU	-	EXPRESSION TAG	UNP P41250
A	687	GLU	-	EXPRESSION TAG	UNP P41250
A	688	HIS	-	EXPRESSION TAG	UNP P41250
A	689	HIS	-	EXPRESSION TAG	UNP P41250
A	690	HIS	-	EXPRESSION TAG	UNP P41250
A	691	HIS	-	EXPRESSION TAG	UNP P41250
A	692	HIS	-	EXPRESSION TAG	UNP P41250
A	693	HIS	-	EXPRESSION TAG	UNP P41250

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

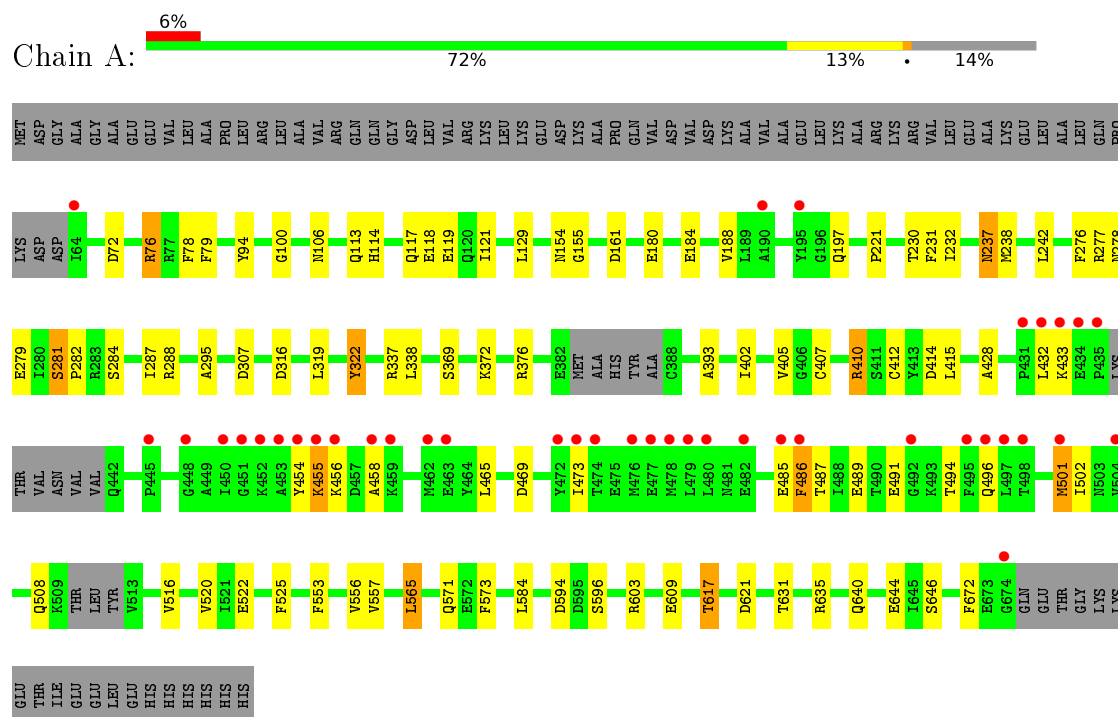
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	100	Total	O	0	0
			100	100		

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.

- Molecule 1: Glycine-tRNA ligase



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	91.24Å 91.24Å 246.61Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.66 – 2.74 34.67 – 2.74	Depositor EDS
% Data completeness (in resolution range)	99.0 (34.66-2.74) 99.0 (34.67-2.74)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.81 (at 2.72Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8_1069)	Depositor
R, R_{free}	0.210 , 0.255 0.209 , 0.253	Depositor DCC
R_{free} test set	1421 reflections (5.07%)	DCC
Wilson B-factor (Å ²)	49.8	Xtriage
Anisotropy	0.086	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 51.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 28044 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4756	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.60% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL

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	0.25	0/4752	0.44	0/6430

There are no bond length outliers.

There are no bond angle outliers.

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	4650	0	4453	48	1
2	A	6	0	8	1	0
3	A	100	0	0	3	1
All	All	4756	0	4461	48	1

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 5.

All (48) 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:617:THR:HG23	1:A:631:THR:HB	1.63	0.79

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:553:PHE:O	1:A:635:ARG:NH2	2.21	0.73
1:A:428:ALA:HB3	1:A:516:VAL:HB	1.75	0.69
1:A:76:ARG:NH2	1:A:609:GLU:OE2	2.32	0.63
1:A:556:VAL:HG13	1:A:557:VAL:HG13	1.81	0.63
1:A:316:ASP:OD1	1:A:337:ARG:NH1	2.32	0.63
1:A:232:ILE:HB	1:A:238:MET:HG2	1.84	0.59
1:A:621:ASP:HB3	2:A:701:GOL:H32	1.84	0.58
1:A:279:GLU:HB3	1:A:282:PRO:HG3	1.84	0.58
1:A:237:ASN:N	1:A:237:ASN:OD1	2.36	0.55
1:A:393:ALA:HB3	1:A:405:VAL:HB	1.89	0.55
1:A:281:SER:OG	1:A:281:SER:O	2.23	0.54
1:A:487:THR:OG1	1:A:496:GLN:OE1	2.22	0.54
1:A:279:GLU:HG2	1:A:288:ARG:HG2	1.90	0.54
1:A:197:GLN:NE2	1:A:221:PRO:O	2.32	0.52
1:A:640:GLN:HB2	1:A:672:PHE:O	2.10	0.52
1:A:154:ASN:OD1	1:A:155:GLY:N	2.44	0.51
1:A:277:ARG:O	1:A:279:GLU:N	2.42	0.51
1:A:106:ASN:HB3	1:A:557:VAL:HA	1.94	0.49
1:A:242:LEU:HB3	1:A:276:PHE:CD1	2.48	0.48
1:A:295:ALA:HB3	1:A:525:PHE:HB2	1.96	0.47
1:A:594:ASP:OD2	1:A:603:ARG:HD2	2.16	0.46
1:A:180:GLU:OE1	1:A:180:GLU:N	2.39	0.46
1:A:410:ARG:NH2	1:A:414:ASP:OD1	2.49	0.45
1:A:410:ARG:NE	3:A:896:HOH:O	2.31	0.45
1:A:72:ASP:O	1:A:76:ARG:HB2	2.17	0.45
1:A:432:LEU:HB3	1:A:433:LYS:H	1.68	0.45
1:A:412:CYS:HB2	1:A:520:VAL:HG23	1.98	0.44
1:A:644:GLU:HG3	1:A:646:SER:H	1.83	0.44
1:A:573:PHE:N	3:A:840:HOH:O	2.50	0.44
1:A:118:GLU:OE2	3:A:841:HOH:O	2.21	0.44
1:A:465:LEU:HD13	1:A:501:MET:HE1	1.98	0.43
1:A:79:PHE:O	1:A:100:GLY:HA3	2.19	0.43
1:A:322:TYR:HB3	1:A:376:ARG:HA	2.01	0.43
1:A:469:ASP:O	1:A:473:ILE:HG13	2.19	0.43
1:A:489:GLU:HB2	1:A:494:THR:HG22	2.02	0.42
1:A:287:ILE:HD13	1:A:402:ILE:HG21	2.01	0.42
1:A:455:LYS:HB2	1:A:456:LYS:H	1.75	0.42
1:A:184:GLU:O	1:A:188:VAL:HG23	2.20	0.41
1:A:412:CYS:SG	1:A:415:LEU:HG	2.60	0.41
1:A:565:LEU:HD21	1:A:603:ARG:HB2	2.03	0.41
1:A:230:THR:OG1	1:A:231:PHE:N	2.54	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:369:SER:HB2	1:A:372:LYS:HE3	2.02	0.41
1:A:119:GLU:HB2	1:A:121:ILE:HG13	2.03	0.41
1:A:113:GLN:HA	1:A:117:GLN:HB2	2.03	0.40
1:A:407:CYS:HA	1:A:522:GLU:O	2.21	0.40
1:A:319:LEU:HD22	1:A:338:LEU:HD13	2.03	0.40
1:A:454:TYR:O	1:A:458:ALA:HB2	2.21	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:94:TYR:O	3:A:850:HOH:O[7_645]	2.16	0.04

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	589/693 (85%)	552 (94%)	27 (5%)	10 (2%)	11 27

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	114	HIS
1	A	455	LYS
1	A	491	GLU
1	A	502	ILE
1	A	284	SER
1	A	571	GLN
1	A	278	ASN
1	A	485	GLU
1	A	486	PHE
1	A	508	GLN

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	490/599 (82%)	475 (97%)	15 (3%)	47 77

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	76	ARG
1	A	78	PHE
1	A	129	LEU
1	A	161	ASP
1	A	237	ASN
1	A	281	SER
1	A	307	ASP
1	A	322	TYR
1	A	410	ARG
1	A	486	PHE
1	A	501	MET
1	A	565	LEU
1	A	584	LEU
1	A	596	SER
1	A	617	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

1 ligand is 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	GOL	A	701	-	5,5,5	0.24	0	5,5,5	0.25	0

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	GOL	A	701	-	-	0/4/4/4	0/0/0/0

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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	701	GOL	1	0

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	597/693 (86%)	0.23	40 (6%) 21 20	26, 50, 107, 129	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	432	LEU	6.2
1	A	450	ILE	5.2
1	A	435	PRO	4.7
1	A	501	MET	4.6
1	A	434	GLU	4.4
1	A	504	VAL	4.3
1	A	451	GLY	4.3
1	A	478	MET	3.8
1	A	476	MET	3.6
1	A	458	ALA	3.5
1	A	455	LYS	3.5
1	A	485	GLU	3.4
1	A	448	GLY	3.2
1	A	454	TYR	3.2
1	A	453	ALA	3.2
1	A	477	GLU	3.2
1	A	479	LEU	3.2
1	A	492	GLY	3.2
1	A	456	LYS	3.1
1	A	190	ALA	3.1
1	A	497	LEU	3.1
1	A	482	GLU	3.1
1	A	495	PHE	3.0
1	A	486	PHE	2.9
1	A	474	THR	2.9
1	A	473	ILE	2.8
1	A	433	LYS	2.8

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	462	MET	2.8
1	A	472	TYR	2.7
1	A	431	PRO	2.6
1	A	452	LYS	2.4
1	A	459	LYS	2.3
1	A	445	PRO	2.3
1	A	480	LEU	2.3
1	A	498	THR	2.2
1	A	674	GLY	2.2
1	A	463	GLU	2.1
1	A	64	ILE	2.1
1	A	496	GLN	2.1
1	A	195	TYR	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 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, 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	LLDF	B-factors(Å ²)	Q<0.9
2	GOL	A	701	6/6	0.86	0.21	2.55	49,56,65,65	0

6.5 Other polymers [i](#)

There are no such residues in this entry.