



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 8WIG / pdb_00008wig
Title : Crystal structure of E. coli ThrS catalytic domain mutant G463S/Q484A
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Deposited on : 2023-09-24
Resolution : 3.22 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
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

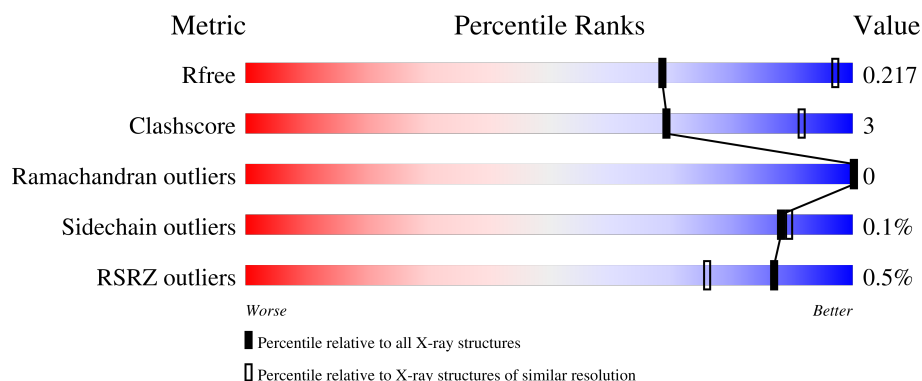
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.22 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	1768 (3.24-3.20)
Clashscore	190562	1879 (3.24-3.20)
Ramachandran outliers	187476	1844 (3.24-3.20)
Sidechain outliers	187428	1843 (3.24-3.20)
RSRZ outliers	180081	1768 (3.24-3.20)

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	410	 89% 9% .
1	B	410	 87% 11% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6539 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 Threonine-tRNA ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	401	Total	C	N	O	S	1	0	0
			3271	2066	575	607	23			
1	B	401	Total	C	N	O	S	0	0	0
			3266	2063	575	605	23			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	241	MET	-	initiating methionine	UNP A0A8S7FUD7
A	463	SER	GLY	engineered mutation	UNP A0A8S7FUD7
A	484	ALA	GLN	engineered mutation	UNP A0A8S7FUD7
A	643	LEU	-	expression tag	UNP A0A8S7FUD7
A	644	GLU	-	expression tag	UNP A0A8S7FUD7
A	645	HIS	-	expression tag	UNP A0A8S7FUD7
A	646	HIS	-	expression tag	UNP A0A8S7FUD7
A	647	HIS	-	expression tag	UNP A0A8S7FUD7
A	648	HIS	-	expression tag	UNP A0A8S7FUD7
A	649	HIS	-	expression tag	UNP A0A8S7FUD7
A	650	HIS	-	expression tag	UNP A0A8S7FUD7
B	241	MET	-	initiating methionine	UNP A0A8S7FUD7
B	463	SER	GLY	engineered mutation	UNP A0A8S7FUD7
B	484	ALA	GLN	engineered mutation	UNP A0A8S7FUD7
B	643	LEU	-	expression tag	UNP A0A8S7FUD7
B	644	GLU	-	expression tag	UNP A0A8S7FUD7
B	645	HIS	-	expression tag	UNP A0A8S7FUD7
B	646	HIS	-	expression tag	UNP A0A8S7FUD7
B	647	HIS	-	expression tag	UNP A0A8S7FUD7
B	648	HIS	-	expression tag	UNP A0A8S7FUD7
B	649	HIS	-	expression tag	UNP A0A8S7FUD7
B	650	HIS	-	expression tag	UNP A0A8S7FUD7

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

- Molecule 1: Threonine–tRNA ligase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	90.13Å 108.18Å 112.42Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	56.21 – 3.22 56.21 – 3.22	Depositor EDS
% Data completeness (in resolution range)	99.6 (56.21-3.22) 99.7 (56.21-3.22)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.24 (at 3.19Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.208 , 0.222 (Not available) , 0.217	Depositor DCC
R_{free} test set	917 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	69.3	Xtriage
Anisotropy	0.536	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 34.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.015 for -h,l,k	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6539	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.65% 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

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	0.08	0/3342	0.24	0/4501
1	B	0.08	0/3337	0.24	0/4494
All	All	0.08	0/6679	0.24	0/8995

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	3271	0	3203	20	0
1	B	3266	0	3198	27	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
All	All	6539	0	6401	45	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 3.

All (45) 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:430:SER:H	1:B:433:MET:HE3	1.56	0.70
1:B:543:VAL:HG23	1:B:590:VAL:HG11	1.78	0.66
1:B:354:ARG:HG2	1:B:386:ILE:HG12	1.80	0.64
1:A:582:ILE:HD13	1:A:609:ARG:HE	1.66	0.60
1:B:252:ASP:OD1	1:B:267:HIS:ND1	2.36	0.58
1:A:538:ALA:O	1:A:568:ARG:NH1	2.40	0.55
1:A:543:VAL:HG23	1:A:590:VAL:HG11	1.90	0.53
1:A:608:VAL:HB	1:A:616:LEU:HB2	1.91	0.52
1:A:343:GLN:HG3	1:B:255:HIS:CE1	2.44	0.52
1:A:296:PRO:HB3	1:B:263:MET:HB3	1.94	0.48
1:B:420:LEU:HD12	1:B:445:LEU:HD12	1.96	0.47
1:A:548:THR:HG23	1:A:550:SER:H	1.78	0.47
1:B:348:TYR:HB3	1:B:499:GLY:HA2	1.97	0.47
1:B:627:LYS:HB3	1:B:640:LEU:HD11	1.97	0.46
1:A:572:ASP:OD2	1:A:585:HIS:NE2	2.43	0.46
1:B:334:CYS:HB2	1:B:335:PRO:HD3	1.97	0.46
1:B:476:ARG:NH2	1:B:527:GLU:OE1	2.42	0.45
1:A:544:ILE:HG12	1:A:594:LEU:HB2	1.97	0.45
1:B:390:GLU:H	1:B:390:GLU:CD	2.25	0.45
1:A:427:ARG:HD2	1:A:434:TRP:CE2	2.51	0.45
1:A:427:ARG:HD2	1:A:434:TRP:NE1	2.32	0.45
1:A:371:HIS:H	1:A:375:ARG:HB2	1.81	0.45
1:B:551:GLN:NE2	1:B:598:ASP:OD1	2.46	0.45
1:B:441:LEU:HB3	1:B:466:ILE:HD11	2.00	0.44
1:B:465:LYS:NZ	1:B:467:GLU:OE2	2.50	0.44
1:A:244:HIS:HB3	1:A:527:GLU:HG3	2.00	0.43
1:B:301:ARG:NH2	1:B:326:GLU:OE1	2.35	0.43
1:B:594:LEU:HD22	1:B:608:VAL:HG22	2.00	0.43
1:A:356:ALA:HA	1:A:384:ALA:HA	2.02	0.42
1:A:428:ILE:HG12	1:A:461:PHE:HD2	1.85	0.41
1:A:533:PHE:O	1:A:611:ARG:NH2	2.44	0.41
1:A:465:LYS:HB3	1:A:484:ALA:HB2	2.01	0.41
1:B:260:ALA:HB1	1:B:263:MET:HE3	2.03	0.41
1:B:374:MET:HA	1:B:520:ARG:HG3	2.02	0.41
1:B:297:PHE:CE2	1:B:360:SER:HB2	2.56	0.41
1:B:304:TRP:CH2	1:B:331:PRO:HD2	2.56	0.41
1:B:394:ARG:HH12	1:B:440:ASP:CG	2.28	0.41
1:A:310:TRP:O	1:A:314:LYS:HE3	2.21	0.41
1:A:535:THR:OG1	1:A:631:GLU:OE2	2.29	0.40
1:B:363:ARG:HE	1:B:363:ARG:HB3	1.77	0.40
1:B:427:ARG:HD2	1:B:434:TRP:CE2	2.56	0.40
1:A:393:ILE:HD11	1:A:508:VAL:HG11	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:317:MET:HG2	1:B:331:PRO:HG3	2.03	0.40
1:B:445:LEU:HD23	1:B:445:LEU:HA	1.97	0.40
1:B:543:VAL:HG11	1:B:585:HIS:CD2	2.55	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	399/410 (97%)	396 (99%)	3 (1%)	0	100	100
1	B	399/410 (97%)	396 (99%)	3 (1%)	0	100	100
All	All	798/820 (97%)	792 (99%)	6 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	355/365 (97%)	355 (100%)	0	100	100
1	B	343/365 (94%)	342 (100%)	1 (0%)	86	87
All	All	698/730 (96%)	697 (100%)	1 (0%)	88	89

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	603	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	268	ASN
1	A	371	HIS
1	A	448	ASN
1	B	333	ASN
1	B	362	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 2 ligands modelled in this entry, 2 are 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 [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	401/410 (97%)	0.08	2 (0%) 87 76	38, 62, 93, 116	98 (24%)
1	B	401/410 (97%)	0.08	2 (0%) 87 76	34, 62, 84, 125	101 (25%)
All	All	802/820 (97%)	0.08	4 (0%) 87 76	34, 62, 88, 125	199 (24%)

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	640	LEU	3.2
1	A	375	ARG	2.4
1	A	297	PHE	2.4
1	B	313	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 oligosaccharides 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	701	1/1	0.99	0.03	58,58,58,58	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	ZN	B	701	1/1	0.99	0.05	62,62,62,62	1

6.5 Other polymers [i](#)

There are no such residues in this entry.