



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 06:19 PM GMT

PDB ID : 1A8H  
Title : METHIONYL-TRNA SYNTHETASE FROM THERMUS THERMOPHILUS  
Authors : Sugiura, I.; Nureki, O.; Ugaji, Y.; Kuwabara, S.; Lober, B.; Giege, R.; Moras, D.; Yokoyama, S.; Konno, M.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)  
Deposited on : 1998-03-26  
Resolution : 2.00 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](mailto: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.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

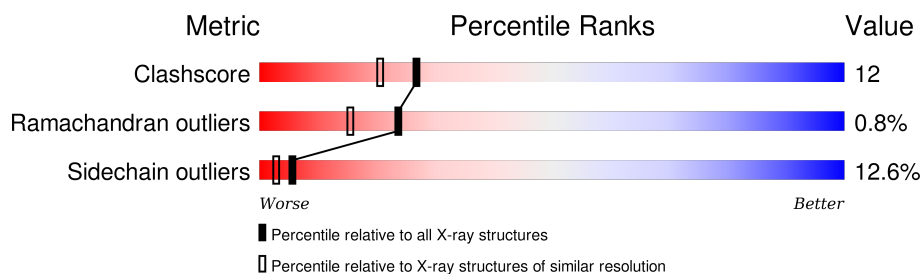
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.00 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))
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)

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  $\geq 3$ , 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	500	 66% 31% •

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4237 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 METHIONYL-TRNA SYNTHETASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	500	Total	C	N	O	S	0	0	0
			4101	2651	708	729	13			

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

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

- Molecule 3 is water.

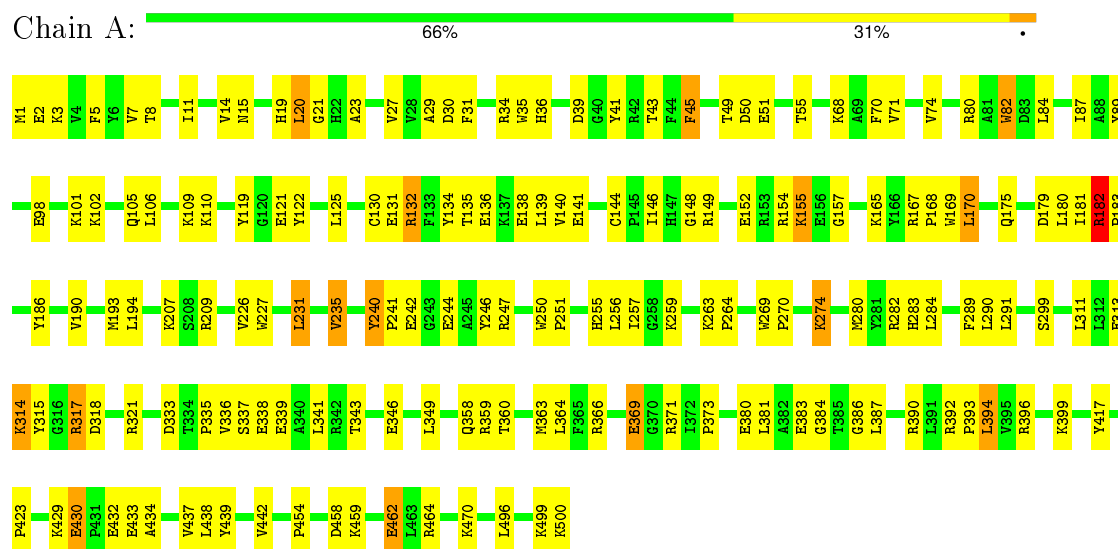
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	135	Total	O	0	0
			135	135		

### 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.

Note EDS was not executed.

#### • Molecule 1: METHIONYL-TRNA SYNTHETASE



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	57.02Å 82.47Å 116.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	6.00 – 2.00	Depositor
% Data completeness (in resolution range)	79.1 (6.00-2.00)	Depositor
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	X-PLOR 3.851	Depositor
R, $R_{free}$	0.205 , 0.283	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4237	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	30.0	wwPDB-VP

## 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	0.30	0/4218	0.54	0/5722

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	4101	0	4059	99	0
2	A	1	0	0	0	0
3	A	135	0	0	1	0
All	All	4237	0	4059	99	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 12.

All (99) 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:235:VAL:HG13	1:A:240:TYR:HB2	1.41	1.02

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:183:PRO:HB2	1:A:186:TYR:HD2	1.36	0.91
1:A:1:MET:HG3	1:A:2:GLU:HG2	1.56	0.87
1:A:182:ARG:HB3	1:A:183:PRO:HD3	1.57	0.84
1:A:384:GLY:HA2	1:A:387:LEU:HD13	1.66	0.77
1:A:125:LEU:HD13	1:A:139:LEU:HD11	1.73	0.71
1:A:349:LEU:HD23	3:A:728:HOH:O	1.93	0.69
1:A:183:PRO:HB2	1:A:186:TYR:CD2	2.24	0.69
1:A:430:GLU:HG3	1:A:433:GLU:HB2	1.75	0.69
1:A:14:VAL:HG11	1:A:51:GLU:HG2	1.73	0.68
1:A:14:VAL:CG1	1:A:51:GLU:HG2	2.23	0.68
1:A:30:ASP:O	1:A:34:ARG:HG3	1.93	0.67
1:A:14:VAL:HG22	1:A:50:ASP:O	1.95	0.67
1:A:386:GLY:O	1:A:390:ARG:HG3	1.96	0.66
1:A:182:ARG:HB3	1:A:183:PRO:CD	2.26	0.65
1:A:19:HIS:HD2	1:A:21:GLY:H	1.43	0.65
1:A:373:PRO:HG2	1:A:439:TYR:HB2	1.79	0.64
1:A:432:GLU:H	1:A:432:GLU:CD	2.02	0.63
1:A:369:GLU:O	1:A:371:ARG:HG3	1.98	0.62
1:A:132:ARG:HD3	1:A:134:TYR:CZ	2.39	0.57
1:A:317:ARG:NH1	1:A:458:ASP:OD2	2.38	0.56
1:A:182:ARG:NH2	1:A:396:ARG:O	2.40	0.54
1:A:181:ILE:HD13	1:A:190:VAL:HG21	1.88	0.54
1:A:146:ILE:C	1:A:148:GLY:H	2.10	0.54
1:A:291:LEU:O	1:A:335:PRO:HA	2.08	0.54
1:A:235:VAL:HG13	1:A:240:TYR:CB	2.26	0.54
1:A:318:ASP:CG	1:A:459:LYS:HG2	2.28	0.54
1:A:98:GLU:OE2	1:A:101:LYS:HD3	2.07	0.54
1:A:130:CYS:O	1:A:131:GLU:HB2	2.08	0.53
1:A:291:LEU:HD11	1:A:333:ASP:HB3	1.90	0.53
1:A:14:VAL:HG13	1:A:49:THR:HB	1.91	0.53
1:A:19:HIS:HD2	1:A:21:GLY:N	2.07	0.52
1:A:257:ILE:HG12	1:A:284:LEU:HD11	1.92	0.51
1:A:314:LYS:HB3	1:A:315:TYR:CD1	2.45	0.51
1:A:144:CYS:SG	1:A:146:ILE:O	2.69	0.50
1:A:360:THR:O	1:A:364:LEU:HG	2.11	0.50
1:A:227:TRP:O	1:A:231:LEU:HB2	2.12	0.50
1:A:263:LYS:HB3	1:A:264:PRO:HD3	1.93	0.50
1:A:259:LYS:HD2	1:A:289:PHE:CE2	2.47	0.50
1:A:45:PHE:CD2	1:A:89:TYR:HA	2.47	0.50
1:A:122:TYR:O	1:A:155:LYS:HA	2.11	0.49
1:A:135:THR:HG23	1:A:138:GLU:OE1	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:167:ARG:HB3	1:A:168:PRO:HD3	1.93	0.49
1:A:8:THR:OG1	1:A:255:HIS:HD2	1.96	0.49
1:A:14:VAL:CG1	1:A:49:THR:HB	2.42	0.49
1:A:70:PHE:O	1:A:74:VAL:HG22	2.12	0.49
1:A:359:ARG:HG2	1:A:363:MET:HE2	1.94	0.49
1:A:179:ASP:O	1:A:283:HIS:HD2	1.96	0.48
1:A:20:LEU:HD13	1:A:290:LEU:HD23	1.94	0.48
1:A:29:ALA:HB2	1:A:256:LEU:HD12	1.95	0.48
1:A:269:TRP:HB3	1:A:270:PRO:HD3	1.95	0.48
1:A:140:VAL:O	1:A:140:VAL:HG23	2.13	0.47
1:A:417:TYR:CE2	1:A:437:VAL:HG13	2.49	0.47
1:A:182:ARG:CB	1:A:183:PRO:HD3	2.37	0.47
1:A:71:VAL:HA	1:A:74:VAL:CG2	2.46	0.46
1:A:263:LYS:N	1:A:264:PRO:CD	2.78	0.46
1:A:146:ILE:O	1:A:148:GLY:N	2.47	0.46
1:A:3:LYS:HD3	1:A:41:TYR:CE1	2.50	0.46
1:A:11:ILE:HG12	1:A:49:THR:C	2.36	0.45
1:A:390:ARG:HH11	1:A:390:ARG:HB3	1.82	0.45
1:A:359:ARG:HG2	1:A:363:MET:CE	2.46	0.45
1:A:101:LYS:O	1:A:105:GLN:HG3	2.16	0.45
1:A:39:ASP:HB3	1:A:41:TYR:CD1	2.52	0.45
1:A:80:ARG:O	1:A:84:LEU:HG	2.17	0.45
1:A:337:SER:OG	1:A:339:GLU:HG2	2.16	0.45
1:A:169:TRP:CE2	1:A:274:LYS:HD3	2.51	0.45
1:A:71:VAL:HA	1:A:74:VAL:HG22	1.98	0.45
1:A:20:LEU:HD11	1:A:336:VAL:HG21	1.99	0.44
1:A:270:PRO:HB3	1:A:280:MET:SD	2.56	0.44
1:A:36:HIS:O	1:A:41:TYR:HB2	2.18	0.44
1:A:11:ILE:O	1:A:11:ILE:HG23	2.17	0.44
1:A:121:GLU:OE1	1:A:155:LYS:HD3	2.17	0.44
1:A:35:TRP:HE3	1:A:454:PRO:HG2	1.83	0.44
1:A:5:PHE:CE2	1:A:7:VAL:HB	2.53	0.44
1:A:240:TYR:HA	1:A:246:TYR:HB2	1.99	0.44
1:A:464:ARG:HD2	1:A:470:LYS:O	2.18	0.43
1:A:119:TYR:OH	1:A:157:GLY:HA3	2.17	0.43
1:A:392:ARG:HB2	1:A:393:PRO:HD3	2.01	0.43
1:A:317:ARG:O	1:A:321:ARG:HG3	2.18	0.43
1:A:31:PHE:CE1	1:A:454:PRO:HB2	2.53	0.43
1:A:423:PRO:HG3	1:A:437:VAL:HG11	2.01	0.43
1:A:341:LEU:C	1:A:341:LEU:HD23	2.39	0.42
1:A:227:TRP:CZ3	1:A:264:PRO:HB3	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:392:ARG:O	1:A:396:ARG:HG2	2.20	0.42
1:A:250:TRP:N	1:A:251:PRO:CD	2.83	0.42
1:A:82:TRP:HA	1:A:82:TRP:HE3	1.85	0.42
1:A:314:LYS:HE3	1:A:338:GLU:OE2	2.20	0.41
1:A:23:ALA:O	1:A:27:VAL:HG23	2.19	0.41
1:A:14:VAL:HG13	1:A:51:GLU:HG2	1.98	0.41
1:A:433:GLU:O	1:A:437:VAL:HG23	2.19	0.41
1:A:394:LEU:HA	1:A:394:LEU:HD12	1.94	0.41
1:A:255:HIS:CE1	1:A:269:TRP:HE1	2.39	0.41
1:A:82:TRP:HA	1:A:82:TRP:CE3	2.56	0.41
1:A:439:TYR:HA	1:A:442:VAL:HG12	2.03	0.41
1:A:82:TRP:HB3	1:A:87:ILE:HB	2.01	0.41
1:A:170:LEU:HD13	1:A:194:LEU:HD11	2.01	0.41
1:A:434:ALA:O	1:A:438:LEU:HG	2.19	0.41
1:A:462:GLU:HB3	1:A:496:LEU:HD11	2.03	0.41
1:A:384:GLY:CA	1:A:387:LEU:HD13	2.43	0.40

There are no symmetry-related clashes.

## 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	498/500 (100%)	466 (94%)	28 (6%)	4 (1%)	24 15

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	226	VAL
1	A	241	PRO
1	A	182	ARG
1	A	240	TYR

### 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	421/421 (100%)	368 (87%)	53 (13%)	<b>5</b> <b>3</b>

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	ASN
1	A	20	LEU
1	A	43	THR
1	A	45	PHE
1	A	55	THR
1	A	68	LYS
1	A	82	TRP
1	A	102	LYS
1	A	106	LEU
1	A	109	LYS
1	A	110	LYS
1	A	132	ARG
1	A	136	GLU
1	A	141	GLU
1	A	149	ARG
1	A	152	GLU
1	A	154	ARG
1	A	155	LYS
1	A	165	LYS
1	A	170	LEU
1	A	175	GLN
1	A	180	LEU
1	A	182	ARG
1	A	193	MET
1	A	207	LYS
1	A	209	ARG
1	A	231	LEU
1	A	235	VAL
1	A	242	GLU
1	A	244	GLU

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Mol	Chain	Res	Type
1	A	247	ARG
1	A	274	LYS
1	A	282	ARG
1	A	299	SER
1	A	311	LEU
1	A	313	GLU
1	A	314	LYS
1	A	317	ARG
1	A	343	THR
1	A	346	GLU
1	A	358	GLN
1	A	366	ARG
1	A	369	GLU
1	A	380	GLU
1	A	381	LEU
1	A	383	GLU
1	A	394	LEU
1	A	399	LYS
1	A	429	LYS
1	A	430	GLU
1	A	462	GLU
1	A	499	LYS
1	A	500	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	ASN
1	A	19	HIS
1	A	61	GLN
1	A	158	ASN
1	A	175	GLN
1	A	221	ASN
1	A	233	ASN
1	A	255	HIS
1	A	283	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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

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 [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](#)

EDS was not executed - this section will therefore be empty.

### 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers [i](#)

EDS was not executed - this section will therefore be empty.